

Fast Reactor Computational Route for Pu Burning Core Design

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**O-arai Engineering Center
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Fast Reactor Calculational Route for Pu Burning Core Design

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Abstract

This document provides a description of a calculational route, used in the Reactor Physics Research Section for sensitivity studies and initial design optimization calculations for fast reactor cores. The main purpose in producing this document was to provide a description of and user guides to the calculational methods, in English, as an aid to any future user of the calculational route who is (like the author) handicapped by a lack of literacy in Japanese. The document also provides for all users a compilation of information on the various parts of the calculational route, all in a single reference. In using the calculational route (to model Pu burning reactors) the author identified a number of areas where an improvement in the modelling of the standard calculational route was warranted - the document includes a description of these changes.

The calculational route makes use of several different computer programs. SLAROM calculates nuclear data from compositions, using either homogeneous or heterogeneous models. CITATION and MOSES do reactor burn-up and/or flux diffusion calculations; CITATION is used for 2D (RZ) calculations, whilst MOSES models 3D (hex-Z) geometry. PENCIL and CITDENS are essentially specialized versions of CITATION (PENCIL includes data preparation and other functions). MASSN calculates fuel cycle mass balances. PERKY performs perturbation and associated calculations, both 1'st order and exact perturbations. JOINT and RZOUT3 provide various dataset interface functions, including energy group condensation.

Briefer descriptions of the calculational route are given, followed by a more detailed step-by-step approach to the calculations. This latter includes examples of all JCL and data files, and a description of all the data that a user may have to employ. The document does not give a complete description of the component programs: where options and/or data are not used in any of the calculations they have generally been ignored; similarly, where data is a fixed value a description has not always been provided.

The document includes comments on and explanations of the modelling assumptions in the various calculations. Practical information on the use of the calculational route and the computer systems is also given.

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プルトニウム燃焼炉心設計ための高速炉解析手順解説

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要旨

この報告書では、炉心技術開発室において用いられている高速炉心の感度解析や初期炉心設計の最適化に関する解析手順について解説する。この報告書の主目的は、将来、この計算手順を追うことになる日本語に不慣れなユーザーの助けとなるように、英語でこの計算手順について解説することにある。また、その他のすべてのユーザーの助けとなるよう、この報告書には、この計算手順の各部分についての詳細な情報も記述した。著者は、プルトニウム燃焼炉心のモデル化のために、基準となる計算手順のモデル化について改良を行った。——これらの変更点についても、本報告書に記載した。——

この解析手順では、複数のコンピュータプログラムを利用する。SLAROM コードは、均質、非均質モデルを用いて、燃料組成から実行断面積を計算する。CITATION コード、MOSES コードは炉心の燃焼方程式や中性子拡散方程式を解く。CITATION コードは、2次元 (RZ モデル) 体系の計算に対して用い、MOSES コードは、3次元 (Hex-Z モデル) 体系計算に用いる。PENCIL コード、CITDENS コードは、CITATION を特殊目的利用のために書き換えたものである。(PENCIL コードは、計算データの準備等の機能を持つ。) MASSN コードは、燃料サイクルの物質収支を計算する。PERKY は、1次摂動計算や厳密摂動計算を行う。JOINT コードと RZOUT3 コードは、中性子エネルギー縮約の機能と計算コード間のデータ交換の機能を持つ。

解析手順について簡単に述べた後、この解析の流れをステップ毎に詳細に説明する。また、付録にはサンプル JCL とデータファイルを添付し、ユーザーが必要なデータについて説明を行った。なお、この報告書は、それぞれの計算コードについては、完全に解説を行っていない。通常、無視できるオプションや、通常、固定値を用いるデータに関する記述は省いた。

この報告書には、種々の計算で行ったモデル化や仮定についてもコメントや説明を記載した。また、この計算手順やコンピュータシステムを利用する上で必要となる実際的な情報についてもあわせて記載した。

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1 **Introduction**

This note describes the calculational route I have been using, after 2 years experience at PNC, to carry out fuel cycle and associated reactor safety parameter calculations. The calculational route is one appropriate to studies of design variations and (initial) design optimization work.

This document was written with two intentions. First, for my own reference and to help me in delineating certain parts of the route which could benefit from the addition or improvement of automated data handling routines. Secondly, as an aid to anyone following me in performing similar calculations and similarly handicapped by a lack of Japanese reading ability. The route as described has been used for calculations on Pu burning cores^(8,9); it is described in such a way as to aid in its adaptation to different fast reactor types.

The note gives three separate descriptions of the calculational route, in increasing detail. First, Section 2 just gives a brief description, as an introduction. Section 3 gives somewhat more detail, concentrating on identifying the relationships of the data files that are transferred between the various component calculations. For a fully detailed description of the calculations see Section 6; included are example JCL and data files (listed in Appendix VII) with instructions for use. Appendices give yet more information, such as brief program manuals for those programs where full manuals only exist in Japanese.

2 A brief description.

As the title suggests, this section gives the simplest description of the calculational route.

The calculations can be considered as split into 6 blocks, not all of which may be required for any particular assessment:

- Iterative calculation to produce condensed nuclear data
- Basic fuel cycle calculation
- Perturbation calculations
- Rod worth calculations
- Shutdown margin calculations
- 3D burn-up calculation.

2.1 The 6 calculation blocks.

The 'iterative condensation' calculations consist of: PENCIL burn-up calculations and SLAROM-JOINT-CITATION (S-J-C) snapshot calculations. The latter uses the 'full' 70-group JENDL-3.2 nuclear data, from which it condenses a new reduced-group dataset, used in the PENCIL and other calculations. The calculations are continued in turn, until Keff converges.

The 'basic fuel cycle' calculations consist of a PENCIL burn-up calculation with reduced groups, then S-J-C snapshot calculations in 70 groups - the latter for BOC & EOC conditions and also for Doppler & Na void conditions. Finally, there is a MASSN mass balance calculation.

The 'perturbation' calculations use PERKY to calculate neutron lifetime and delayed neutron fraction; it can also be used to calculate the contributions to Na void (by reaction) and to Doppler (by nuclide). It first requires one SLAROM and several CITATION calculations to prepare data of the required form.

'Rod worth' calculations use the 3D model MOSES in steady-state mode. Reduced-group nuclear data are used. Keff values are calculated for a variety of rod positions.

Various factors, in addition to the foregoing rod worth calculations, are needed for 'shutdown margin' calculations. S-J-C calculations with mesh sizes fractionally increased give reactivity factors for core size expansion. The temperature effect on nuclear data can be obtained from either an S-J-C calculation as part of the 'basic fuel cycle', or from a MOSES 'rod worth' calculation; whichever is used must employ nuclear data for cold shutdown temperatures. Material density reactivity coefficients are obtained from a special package based on PERKY; it uses one SLAROM and 2 CITATION calculations to set up required data.

For accurate values of peak pin rating, the PENCIL burn-up calculation has to be repeated using a MOSES '3D burn-up' calculation. Even with a reduced number of energy groups, this calculation takes a long time to run.

2.2 The computer programs.

MOSES & CITATION are diffusion models, allowing both steady state and burn-up calculations. MOSES is a 3D (hex-Z) geometry: it allows a 6-mesh per S/A model, but I am only using it in a 1 mesh per S/A mode, with interpolation of rating (but not number densities) to 6 meshes per S/A (based on the Askew method). CITATION does support 3D calculations, but is only being used in 2D (RZ) geometry here. [As an old code, I guess it is written to use disk storage more than is necessary with modern computers, so large models take a long time (elapsed time much more so than CPU time).]

SLAROM is the cell model, used to adjust the infinite dilution cross-section data of JENDL-3.2 (the basic 70-group data used in all the calculations) to the compositions of the particular model zones. It can model heterogeneous as well as homogeneous cells.

JOINT is a set of editing programs used as steps of batch jobs. It also does such things as condensing the nuclear data to a reduced number of energy groups, and altering nuclear data between different storage formats.

PENCIL is a program which basically comprises running a CITATION burn-up calculation, which models a series of refuelling steps until fuel cycle equilibrium is achieved. What PENCIL also does is to adjust Pu enrichments and then repeat the CITATION calculation, iterating on Pu enrichments until target Keff and rating balances (inner to outer zone peaks) are achieved.

CITDENS is a slightly modified form of CITATION: it is used to repeat the final CITATION calculation from a PENCIL run, producing an output file of number densities at each step throughout the calculation.

MASSN is basically an editing program; it takes the number density data produced by CITDENS and transforms it into information on the mass balances during the fuel cycle.

PERKY does perturbation calculations - both exact and 1'st order - and also does the associated calculations of neutron lifetime and delayed neutron fraction. The special PERKY package to produce density reactivity coefficients is basically a program script for linking a PERKY calculation with a whole series of data files, and incorporating pre- and post-processors and a CITATION calculation.

RZOUT3 is a short editing program, run interactively to

extract number density data for a single snapshot condition from the output file created by CITDENS.

There are various places where a significant amount of hand editing has to be done to transform the output files produced by one calculation into input required by another. I am in the process of trying to automate these, where practicable, see Section 7. (This note is in part being written to help me define the requirements.)

2.3 Some general comments.

In the most recent series of calculations that I undertook, I significantly increased the detail in which things were modelled, compared with both my earlier calculations and with the standard route as described in PNC documentation (PN9520 94-003). Some of these changes were made because earlier results had pointed up weaknesses in the calculations; others were changed because the nature of the study now called for more detailed but less wide ranging calculations. I will try to make clear where there is a reasonable alternative to use more or less detailed calculations, depending on the aims of the particular study being carried out.

Less detailed calculations can model just half core height (with a reflective centre plane), including no rod absorber in the model. This option was used in my earlier calculations (it is part of the standard PNC route); it is fine for parametric survey type work, but introduces systematic biases into any definitive calculations. The route as I describe it here models the full core height and includes absorbers: using the simpler model will require an extra step to produce nuclear data for the 3D MOSES calculations, in order to incorporate data for absorber material.

The standard route used 7-group data for PENCIL/CITDENS burn-up calculations; this introduced significant random uncertainties in Keff (> half a percent). I am now using 18 groups for these burn-up calculations. The effect may well have been largely a consequence of the core designs I am examining (moderator in the core, no breeder zones). If the use of 18-group data makes the PENCIL run times excessive, then it is always possible to run PENCIL twice, first a preliminary calculation with 7-group data, to get Pu enrichments to use as starting values in the definitive PENCIL calculation (with 18 groups). This should minimize the calculation time.

The level of detail in the flux and number density meshes will to some extent depend on the type of core design. The values used were for a breederless core, where edge leakage effects are particularly important, hence the refined core edge meshes.

3 The calculational route - an outline

In the following, I give a more detailed list of the calculations involved in each of the 6 identified blocks, including those input/output files used to connect the various component calculations. Later on I will endeavour to give a full description, including example data and JCL files. The following section is intended to give a view on how the various components fit together.

3.1 Iterative Condensation block

Produces both 7-group and 18-group condensed nuclear data. If the effects of temperature change on nuclear data (part of the shutdown margin calculation) are to be calculated using MOSES rather than using S-J-C [and I recommend that it is, though temperature reactivity values for the 2 methods differ by $<0.1\% \Delta k/k'$], then a 3'rd condensed nuclear dataset is produced, in 7 energy groups, but for shutdown temperatures.

The following series of calculations cycles around repeatedly, until there is little change in K_{eff} (either PENCIL or CITATION) between successive calculations. Only on the final iteration does the S-J-C calculation need to produce all 3 condensed nuclear data files.

PENCIL burn-up calculations in 18 groups, (seeks optimum Pu enrichment)

input	18-group nuclear data 'PENCIL data' file 'CITATION data' file 1 (burn-up)
output	temporary file to run CITDENS

CITDENS burn-up calculation (repeats final case from PENCIL)

input	temporary file from PENCIL
output	number density file (multiple entries)

RZOUT3 extracts single entry (BOC) from CITDENS output

Hand Edit transform RZOUT3 output to 'SLAROM data' format

S-J-C: SLAROM produces effective cross-sections (70 groups); CITATION snapshot calculation in 70 groups; JOINT condenses nuclear data (using flux from CITATION)

input	JENDL-3.2 70 group nuclear data 'SLAROM data' from Hand Edit 'CITATION data' file 2 (snapshot) 'JOINT data' flux extraction file 'JOINT data' condensation file 1 (7 group) 'JOINT data' condensation file 2 (18 group)
-------	--

```

                                'JOINT data' reformatting file 1
                                                (7 group)
                                'JOINT data' reformatting file 2
                                                (18 group)
output    temporary flux file, 70 groups
          PDS format nuclear data, 7 groups
          PDS format nuclear data, 18 groups
          'micro' nuclear data, 7 groups
          'micro' nuclear data, 18 groups

Hand Edit    alter 'SLAROM data' to cold
                temperatures

S-J-C: as previous case, except is for cold
temperatures
input      JENDL-3.2 70 group nuclear data
          'SLAROM data' for cold temperatures
          'CITATION data' file 2 (snapshot)
          'JOINT data' flux extraction file
          'JOINT data' condensation file 1
                                                (7 group)
          'JOINT data' reformatting file 1
                                                (7 group)
output    temporary flux file, 70 groups
          PDS format nuclear data, 7 group, cold
          'micro' nuclear data, 7 group, cold
    
```

The "'micro' nuclear data, 18 groups" output, is used as the 18 group input file for PENCIL in the next iteration. The original 18 group input file, and the 70 group JENDL-3.2 input, are pre-existing nuclear data files. All other input files have to be created at the start of the calculation, but are invariant between iterations.

Note that whilst on the 1'st iteration the PENCIL calculation may be allowed to find optimum Pu enrichments, on all subsequent iterations the enrichments must be fixed at those used in the first CITDENS calculation.

The output files generally need to be given new names with each iteration, to avoid conflicts.

It is not necessary to go through this process of condensing new nuclear data for every core variant examined: the main indicators of a requirement for new data are if there are any different nuclides present, or if there are large changes in the proportions of those present. An obvious test for whether a new condensation is necessary: use an 18-group nuclear dataset as the starting point in the first iteration, if the 2'nd iteration PENCIL doesn't change Keff much then there is no need to recondense (since you will already have the slightly improved data produced by the 1'st iteration, this approach is only really useful to check one case that is representative of several others).

3.2 Basic Fuel Cycle block

When the above iterative process has converged on final condensed nuclear datasets - or, if a pre-existing condensed dataset is considered to be adequate for the core/fuel cycle variant being modelled - the following series of calculations is performed.

PENCIL burn-up calculations in 18 groups, (seeks optimum Pu enrichment)

input	18-group nuclear data 'PENCIL data' file
output	'CITATION data' file 1 (burn-up) temporary file to run CITDENS

CITDENS burn-up calculation (repeats final case from PENCIL)

input	temporary file from PENCIL
output	number density file (multiple entries)

MASSN calculation of fuel cycle mass balances

input	number density file from CITDENS
-------	----------------------------------

RZOUT3 extracts single entries (both BOC & EOC) from CITDENS number density file

Hand Edit transform RZOUT3 output to 'SLAROM data' format; 4 or 5 separate files created:

	BOC, unmodified
	EOC, unmodified
	EOC, Na void (0.0 Na number density)
	EOC, Doppler (fuel temperature +500)
(optional)	BOC, cold shutdown (all temperatures 473K)

! S-J-C: SLAROM produces effective cross-sections (70 groups); CITATION snapshot calculation in 70 groups;
! JOINT is just the interface between the 2 other programs
!
! input JENDL-3.2 70 group nuclear data
! 'SLAROM data' from Hand Edit
! 'CITATION data' file 2 (snapshot)

An S-J-C calculation is done for each 'SLAROM data' file created by the Hand Edit.

The cold shutdown case is done if the effects on nuclear data of the temperature change are being calculated by CITATION: if MOSES is being used, cold 7-group nuclear data will have been condensed in the relevant 'iterative calculation', so no cold shutdown S-J-C calculation need be done here.

All the data files not created within this block are identical to the equivalent data files in the 'iterative condensation' block. The JENDL-3.2 and 18-group nuclear data files are pre-existing (the 18-group file comes from an 'iterative condensation' calculation). The one PENCIL and two CITATION input files have to be created by hand for the specific reactor and fuel cycle variant.

3.3 Perturbation block

SLAROM and CITATION calculations produce flux and PDS format nuclear data for the 'normal' reactor state and for the perturbed states ('Na void' and 'Doppler' conditions), all for EOC. An exact perturbation calculation using the full 70 group nuclear data is required for the Na void worth to be calculated accurately (compared to a difference in Keff from S-J-C). The Doppler perturbation remains accurate with a 1'st order perturbation calculation in 18 energy groups. The standard PNC delayed neutron data is in 18 energy groups, hence the use of 18 groups in the delayed neutron fraction calculation.

RZOUT3 extracts a single entry from number density file
input CITDENS n.d. output of 'basic fuel cycle'
output number density data for EOC

Hand Edit transform RZOUT3 output to 'SLAROM data' format, including versions for 'normal', 'Doppler' and 'void' core states

SLAROM produces 70 group effective cross-sections (for 'normal', 'Doppler' and 'void' states)

input JENDL-3.2 70 group nuclear data
'SLAROM data' from Hand Edit
output 70 group nuclear data file ('normal', 'Doppler' and 'void')

CITATION, direct and adjoint flux snapshot calculations, for the 'normal' core state and in 70 energy groups, plus associated JOINT condensation to 18 energy groups

input 70 group nuclear data from SLAROM
'CITATION data' file, 70 groups, 'normal'
'JOINT data', flux extraction file, 'normal'
'JOINT data', condensation file, 'normal'
output temporary flux file (70 groups)
direct & adjoint flux file, 70 groups, 'normal', (FT34 format)

PDS format nuclear data, 18 groups,
 'normal'

CITATION, direct flux snapshot calculation, for the
 'void' core state and in 70 groups
 input 70 group nuclear data from SLAROM
 'CITATION data' file, 'void'
 output direct flux file, 70 groups, 'void',
 (FT34 format)

CITATION, direct and adjoint flux snapshot
 calculations, for the 'normal' core state and in 18
 energy groups
 input 18 group 'normal' PDS data - from 70
 group 'normal' CITATION calc.
 'CITATION data' file, 18 groups,
 'normal'
 output direct & adjoint flux file, 18 groups,
 'normal', (FT34 format)

CITATION, direct flux snapshot calculations, for the
 'Doppler' core state and in 70 energy groups, plus
 associated JOINT condensation to 18 energy groups
 input 70 group nuclear data from SLAROM
 'CITATION data' file, 70 groups
 'Doppler'
 'JOINT data', flux extraction file,
 'Doppler'
 'JOINT data', condensation file,
 'Doppler'
 output temporary flux file (70 groups)
 PDS format nuclear data, 18 groups,
 'Doppler'

PERKY prompt neutron lifetime calculation
 input 70 group nuclear data from SLAROM
 'PERKY data' file (neutron lifetime)
 direct & adjoint flux file, 70 groups,
 'normal'- from CITATION

PERKY delayed neutron fraction calculation
 input PDS data file, 18 group, 'normal' -
 from CITATION
 'PERKY data' file (delayed neutron)
 direct & adjoint flux file, 18 groups,
 'normal' - from CITATION

PERKY Na void exact perturbation calculation
 input 70 group nuclear data from SLAROM
 'PERKY data' file (void perturbation)
 direct & adjoint flux file, 70 groups,
 'normal' - from CITATION
 direct flux file, 70 groups, 'void' -
 from CITATION

PERKY Doppler 1'st order perturbation calculation, by
 zone

```

input      PDS data file, 18 group, 'normal' -
           from CITATION
           PDS data file, 18 group, 'Doppler' -
           from CITATION
           'PERKY' data file (Doppler, by zone)
           direct & adjoint flux file, 18 groups,
           'normal' - from CITATION

PERKY Doppler 1'st order perturbation calculation, by
isotope
input      PDS data file, 18 group, 'normal' -
           from CITATION
           PDS data file, 18 group, 'Doppler' -
           from CITATION
           'PERKY' data file (Doppler, by
           isotope)
           direct & adjoint flux file, 18 group,
           'normal' - from CITATION
    
```

The calculations of Doppler contributions are not really necessary, now. The original PNC method allowed only a single temperature for the whole fuel cell, so Doppler effect was overestimated by incorporating effects from steel nuclides. I am now using a heterogeneous 2-region cell model for the fuel, with the fuel pellet separate from other materials, and Doppler temperature changes restricted to this fuel region. Therefore, the effect which the 'by nuclide' analysis was there to examine, has been eliminated. Unless you decide you need to do such a calculation, then several of the above steps can be eliminated: the SLAROM calculation need not include materials for the perturbed Doppler condition; the 70-group CITATION calculation and condensation to 18 groups for the Doppler state can be eliminated, as can both PERKY Doppler perturbation calculations.

3.4 Rod Worth block

For a new core/rods configuration, it will be necessary to survey the range of PCR (main rod) and BCR (backup rod) positions, to find the most effective positions for single stuck rods.

Shutdown margin analysis is restricted to assessing the worths of rod groups with a single rod stuck (there is no consideration of faults during or as a result of rod withdrawal for maintenance). As a result just the following rod positions are to be assessed:

PCRs fully out	BCRs fully out	
PCRs 50% in	BCRs fully out	
PCRs fully in	BCRs fully out	1 PCR fully out
PCRs 50% in	BCRs fully in	1 BCR fully out

The calculations are done for BOC conditions (when shutdown margins are most difficult to meet). If the effect of cold shutdown temperatures on nuclear data is being calculated

using MOSES, the second case is repeated but using the nuclear data for the cold temperature.

PDSDUMP, program to read data - in this case fission spectrum - from PDS format files.

```

input      JCL file contains data
           PDS format nuclear data file, 18 group
           from 'iterative convergence'
output     fission spectrum data, 18 groups
    
```

RZOUT3 extracts single entry from a number density file

```

input      CITDENS n.d. output of 'basic fuel
           cycle'
output     number density data for BOC
    
```

Hand Edit transform RZOUT3 output to 'MOSES data' number density format; condense fission spectrum to 7 groups; use to create 'MOSES data' snapshot file

```

! MOSES snapshot calculation in 7 groups
!   input      'micro' nuclear data, 7 groups (from
!               'iterative condensation')
!   output     'MOSES data' snapshot file from Hand
!               Edit
    
```

A MOSES calculation is done for each of the 4/5 rod configurations and temperature conditions described earlier.

The only differences between the 4 'MOSES data' input files is a (simple) re-positioning of rods; the basic dataset does have to be created by hand for each reactor/fuel cycle. The other input files all come from earlier calculation blocks, as noted.

Rod worths are calculated from the difference in Keff values of the appropriate MOSES calculations.

3.5 Shutdown Margin block

The effect of cold shutdown temperatures on nuclear data is dealt with in other calculation blocks. Two sets of calculations are dealt with here: reactivity coefficients for changing core size; material density reactivity coefficients.

First the core size reactivity coefficient calculations:

```

! Hand Edit      alter radial mesh size (increase 1%)
!   input      'CITATION data' file, snapshot data
!               (file 2) from 'basic fuel cycle'
!   output     'CITATION data' file, +1% radial mesh
    
```

```

!   S-J-C: SLAROM produces effective cross-sections (70
!   groups); CITATION snapshot calculation in 70 groups;
!   JOINT is just the interface between the 2 other
!   programs
!       input      JENDL-3.2 70 group nuclear data
!                  'SLAROM data' file, BOC, from 'basic
!                  fuel cycle'
!                  'CITATION data' file, +1% radial mesh,
!                  from Hand Edit
!

```

The above is repeated, but with the +1% change applied to axial rather than radial meshes.

The remainder of this block is the material density reactivity coefficient calculation:

```

RZOUT3 extracts single entry from a number density
                                     file
input      CITDENS n.d. output of 'basic fuel
                                     cycle'
output     number density data for BOC

```

```

Hand Edit alters RZOUT3 output to 'SLAROM data'
format, includes 'Doppler' version of each material
zone

```

```

SLAROM creates effective cross-sections, in 70 groups
input      JENDL-3.2 70 group nuclear data
           'SLAROM data' from Hand Edit
output     70-group nuclear data file ('normal' &
           'Doppler')

```

```

CITATION calculation for 'normal' core state, plus
JOINT condensation to 18 groups
input      70-group nuclear data file from SLAROM
           'CITATION data', snapshot file,
           'normal'
           'JOINT data', flux extraction file
           'JOINT data', condensation file
           'JOINT data', reformatting file
output     temporary flux file, 70 groups
           PDS format nuclear data, 18 groups,
           'normal'
           'micro' nuclear data, 18 groups,
           'normal'

```

```

Hand Edit, changing zone names from 'normal' to
'Doppler' versions.

```

```

Edit Files:  'CITATION data', snapshot file
             'JOINT data', condensation file

```

```

CITATION calculation for 'Doppler' core state, plus
JOINT condensation to 18 groups
input      70-group nuclear data file from SLAROM
           'CITATION data', snapshot file
           ('Doppler' version, from Hand Edit)
           'JOINT data', flux extraction file

```

'JOINT data', condensation file
('Doppler' version, from Hand Edit)
output temporary flux file, 70 groups
PDS format nuclear data, 18 groups,
'Doppler'

PDSDUMP, program to read data - in this case fission spectrum - from PDS format files.

input JCL file contains data
PDS format nuclear data file, 18 group
'normal', from CITATION
output fission spectrum data, 18 groups

PERKY package, including pre- and post- processors, and a CITATION direct and adjoint flux calculation

input CITDENS n.d. output of 'basic fuel cycle'
Temporary file from PENCIL, used in 'basic fuel cycle' CITDENS
'Fission Spectrum data' file, from PDSDUMP output
PDS format nuclear data, 18 groups, from 'normal' CITATION
PDS format nuclear data, 18 groups, from 'Doppler' CITATION
'micro' nuclear data, 18 groups, from 'normal' CITATION
'Delayed neutron data' file
'Nuclide ID data' file
'Material zone data', 2 files
'Post-processor control data', 3 files
output Several temporary files created by the package (delete them)

The SLAROM run has each zone represented twice, with the 2'nd (Doppler) version having a different zone name - only for fuel/breeder zones is there a temperature difference between the 2 versions. The 'CITATION data' file and 'JOINT data' files used in the normal CITATION calculation were copied directly from the 'iterative condensation' block - if that step was omitted, then these files have to be created now for the reactor/fuel cycle under assessment. The data in the PDSDUMP JCL file is reactor design dependent. The 'delayed neutron data' file is a pre-existing (binary) file.

3.6 3D Burn-up Calculation block

The insertions of different rings of PCRs (main rods) can be used to balance the inner:outer peak ratings; also, different BOC & EOC rod insertions can be used to give a suitable Keff (possibly 1.0) in both states (the PNC reactivity target is 1.0048 at EOC, all rods fully out). It is possible to first use CITATION, rather than the slower running MOSES calculations, to find approximate

values for rod insertions.

```
!   S-J-C: essentially a repeat of a 70-group calculation
!   from the 'basic fuel cycle'
!       input      JENDL-3.2 70 group nuclear data
!                   'SLAROM' data from 'basic fuel cycle'
!                   'CITATION' snapshot data file with
!                               modified rod insertions
```

The above calculations are done for both BOC and EOC states; cases are repeated as necessary, with adjustments, to find acceptable rod positions. These can only give approximate rod insertions, because of feedback effects from altering rod insertions during burn-up.

```
!   Hand Edit, to produce 'MOSES data' burn-up file
!       input      'MOSES data' snapshot file, from 'rod
!                               worth' calculation
!                   Clean core number density data, taken
!                   from 'CITATION' data used in PENCIL
!                   Burn-up modelling information
!       Output     'MOSES data' burn-up file
!
!   MOSES burn-up calculation, basically the same
!   calculation as PENCIL, but in 3D
!       input      'MOSES data' burn-up file from Hand
!                               Edit
!                   'micro' nuclear data, 7 groups, from
!                               'iterative condensation')
```

The above is repeated as necessary, to fine-tune rod insertions (Keff should not be fine-tuned: the 7-group MOSES calculations will introduce a significant Keff error c.f. the 70-group CITATION calculations).

3.7 Additional Calculation block

As was noted earlier, should the 'iterative condensation' and 'basic fuel cycle' blocks be carried out for a half-height core model with no rod absorber present, then an additional calculation block will be necessary to provide absorber nuclear data for any 3D MOSES calculation, whether 'rod worth' or '3D burn-up'. Such a calculation is described here; note that there is no example dataset nor any detailed description in Section 6.

The calculation uses as input the 7-group PDS format nuclear data files created in the final S-J-C calculation of the 'iterative condensation' block; if files need re-creating, then the S-J-C calculation is re-run. To create condensed nuclear data files, in the binary file format used by PENCIL/CITATION and including the previously absent rod absorber material, the S-J-C calculation is repeated, but with certain changes:

The SLAROM calculation includes an extra material - the rod absorber, it is modelled homogeneously (if '3D burn-up' calculations are required, both homogeneous and heterogeneous versions of the rod absorber are modelled).

In CITATION a new material zone is added: a central region, of radius equivalent to 1 S/A, has the material replaced by rod absorber (the homogeneous version).

The JOINT1 step extracts flux data, just as before. Since data is for MOSES, only the 7-group JOINT2 condensation step is required. Only for the rod absorber region is the condensation carried out (twice, if the SLAROM step contained both homogeneous and heterogeneous data), the flux from the new central absorber zone is used for the condensation. Rather than create a new PDS file, the condensed data is added to the pre-existing PDS file.

The MICRO step, transforming PDS format nuclear data to the binary file format used by PENCIL/CITATION, does so for every zone with data now in the PDS file.

The calculation is done twice, for both the normal and cold shutdown temperature conditions.

4 The Computer System.

Batch jobs are run on the Oarai FACOM mainframe computer. There are a number of idiosyncrasies of the FACOM computer system - they can no doubt all be laid at the door of the IBM machines that it copies - which those who have become used to UNIX type systems will have to (re)learn.

The main problems are to do with file storage: not only is the amount of disk space available very restricted (of order a hundred Mb, at best), it is necessary for the user to do a lot of the housekeeping work. Library files (i.e. directories) are defined as a set size when you create them - actually a base size plus expansion block size, but there is a limit of 15 expansion blocks - none of this can be re-set. Once a library has been allocated, and used, a certain amount of space, then that space is permanently assigned to that library (unless the library is deleted) and not available to any other, even if files in the library are deleted. Libraries require regular cleaning, otherwise they will eventually become full - this is not just a question of deleting unwanted files, to release the space it is necessary to run a compression routine. It is necessary to make sure that a library has a suitable setting for record length and file organization for the type of files it will contain.

There is a limit (~15) on the number of batch jobs any uid can have in the queues - this includes output from jobs that have already run.

Sometimes, when a job/process is to create an output file, it is helpful/necessary to have a file of that name already existing, to define the format of the output file.

Error messages are the usual IBM gobbledygook - there was a (Japanese) manual somewhere in the next room which translates these messages, but I have not been able to find it recently. A Japanese colleague may well be able to recognize an error message from the error number given.

With the limited storage for permanent files, it is often necessary to use the temporary file space that is allocated to the section. This, even more than the permanent space, is subject to competition, with multiple users trying to use the same space concurrently - it can be necessary to co-ordinate your calculations so as not to clash for file space with other section members. The temporary filespace is called 'weekly' storage, but if it is full it may be cleared out more frequently by the operators (the same applies to batch job output - it is kept on-line for a nominal week, but it can sometimes be deleted earlier).

Be aware that you should do your own archiving of files kept on the FACOM. There is a tape cassette reader in the computer building to which you have direct access. For storage of batch job output, I now transfer this (FTP) to the Macintosh on my desk (in 2 steps, via the section's

Unix workstation), and store the files on 230 Mb demountable MO disks. Whilst files can be sent back to the FACOM, I have experienced some problems with this, so would not suggest using this route to archive data/JCL files that might want retrieving to the FACOM.

5 Comments on the Calculations.

There are essentially two separate sets of programs for running the calculations. The hallmark of which system is being used is the number of characters in the names used to identify the different material zones of the reactor model, either 5 or 6 characters.

The older versions of the programs use 6-character names: these programs allowed only a single character to identify different isotopes within cross-section data, as a result a maximum of 34 isotopes were allowed. The later version re-assigns one character, using 5 to identify zone and 2 for the isotope - a maximum of 99 isotopes is now possible. The larger number of isotopes is particularly necessary for calculations dealing with minor actinide burning. Note that if a zone name in an input file is less than the required 5/6 character, the programs fill the blanks with '\$' signs; also, excess characters may be truncated; so the number of characters in data in an input file is not an infallible guide to the program version being used. (Note that on some Japanese machines, 'dollar' and 'yen' symbols are interchanged, on others this applies to the 'yen' and 'backslash'.)

Not all programs are available in the updated (5-character) format. Both for the purposes of changing between the 5 and 6-character calculations, and to adopt any updated/corrected programs that become available, it is generally sufficient to just change the program module names in the JCL files.

The 70-group nuclear data is a single set of data which is used for all material compositions. When condensed (7 or 18 group) datasets are produced, these include separate data entries for a number of different material compositions (i.e. each condensed from 70 groups using a flux from a zone with the appropriate composition). In (some) condensed datasets, the data for each material zone include microscopic cross-section data for all 150 nuclides of the JENDL library; because of the bulky nature of this data, such condensed datasets can only represent a limited number of different material zones. In particular, these datasets only have one entry for each fuel Pu enrichment zone, despite the variation in composition across such a zone (a consequence of the burn-up distribution).

Because of the above nature of the condensed datasets, the 'iterative condensation' S-J-C calculations (used to produce the condensed datasets) used a simple geometric representation, with the number of material zones corresponding to the number of condensed cross-section dataset entries - thus each core enrichment zone has a uniform burn-up. The remainder of the CITATION based calculations use a more detailed material zone representation, with axial and radial burn-up (i.e. number density) variation modelled. I found that a larger level of detail than had previously used in the PNC route was

necessary: this was because of the importance of leakage (especially in Na void) to the breederless cores I was modelling - it was sensitive to the number density variation near the core edge.

In the original PNC calculational route (which used fewer material/number density zones to represent the core), the calculations of data for the PERKY perturbations included the production of condensed cross-section data. With the increased number of zones in my models, this encountered problems over trying to create files large enough to include the full 150 nuclide microscopic data for each material zone. I modified the calculation slightly, so that microscopic cross-section data were only stored for a single material zone (one for each of the normal, Doppler and Na void core states), these values being used for all zones in the PERKY calculations.

In the MOSES burn-up calculations, the radial burn-up/number density mesh is automatically set at 1 mesh per S/A by the program. The axial number density mesh corresponds to the AXIAL input parameter (8 meshes in the example datasets). Thus, the MOSES burn-up calculation has a much more detailed material mesh representation than the PENCIL and S-J-C (RZ) calculations, as well as a more accurate flux/rating representation.

In the rod worth (snapshot) MOSES calculations, I originally used a restricted total of just 15 zones for all materials, averaging data from the more detailed CITATION calculations to provide number density input. This restriction was due to a misunderstanding of how MOSES worked; the program is not in fact limited to 15 zones of number density input. I would recommend as far as possible maintaining the same zones as are used in the CITATION calculations. It is only practical to consider radial n.d. zones in MOSES that consist of S/A rings (or part rings, where caused by material zone boundaries). To transfer number densities from CITATION to MOSES, some of the narrower CITATION radial zones may have to be combined.

The radial (number density) mesh boundaries used in CITATION (RZ geometry) should be carefully chosen to correspond (where possible) to S/A ring boundaries in the hex-Z geometry of MOSES.

The 'perturbation' block of calculations is carried out using the original, 6-character, versions of the programs - a version capable of handling large arrays (i.e. a large model) was needed, and some 5-character program versions could not meet the requirements. These perturbation calculations are also limited to a maximum of 40 material zones in the model (the limit comes from some of the JOINT programs), it was necessary to adopt a simplified version of the zone structure used in the 'basic fuel cycle' and other calculations - it was still significantly more detailed than that used in the 'iterative condensation' calculations. (The PERKY package-based calculations for

the 'shutdown margin' block make use of the newer, 5-character, name versions of the programs; it also uses the fully detailed material zone model.)

The program restriction on numbers of material zones that is most likely to be limiting comes from PENCIL: no more than 24 material zones in either Pu enrichment zone.

The example datasets have both fuel and absorber zones modelled heterogeneously in the SLAROM cell lattice calculations. The fuel heterogeneous model is included so that non-fuel materials can be excluded from the temperature changes of Doppler calculations; the more accurate modelling of fuel self-shielding is an added bonus (the effect is minor). The original PNC calculational route included no method for absorber rod homogenization (not a problem in those calculations, since absorber was only modelled in the 3D MOSES rod worth calculations, and one of the rod worth adjustment factors applied was for rod homogenization). I considered it important to include the absorber in the more detailed model I was using, especially since I was looking at B₄C diluent in the core, and its effectiveness could be affected by the presence of the same material in the rods. The heterogeneous cell model used for absorber was designed for fuel element modelling, however, it appears to give reasonable results, and it is important to have some method of representing absorber heterogeneity effects.

I have continued to use homogeneous absorber in the MOSES rod worth calculations (not the burn-up), rather than alter the rod worth adjustments to remove the compensation for absorber heterogeneity effects. (It might be a good idea to change this, just using heterogeneously modelled rod absorber throughout.) The method I have used requires two sets of absorber nuclear data to be included in each condensed nuclear dataset: the S-J-C calculation of the 'iterative condensation' just models heterogeneous absorber in the CITATION calculation, both heterogeneous and homogeneous absorber models are included in the SLAROM calculation, and both are condensed by JOINT, using the single absorber flux calculated by CITATION.

The MOSES calculations have generally used the option for 1 mesh per S/A, with the ratings interpolated to 6 meshes per S/A (using the Askew method). The more detailed 6 mesh per S/A option works, but the extra computer time used does not much alter the results obtained. In some cases I had to use the more detailed calculation, because the other option would not converge. The MOSES option to include an interpolation of number density as well as flux to 6 meshes per S/A is not yet implemented in the program.

As to the insertion of the control rods into the core, for certain calculations these are fixed. The 3D MOSES 'rod worth' cases have rod positions defined as a key part of the calculations. For the PENCIL calculations, rods must be fully withdrawn, since the reactivity target is defined

for that condition. In the case of the '3D burn-up' MOSES calculations, rod insertions are to be adjusted to give the desired Keff value and radial rating profile. As to the remaining calculations - those involving CITATION - I have adopted the following rod insertions: fully withdrawn for cases used in reactivity difference calculations (this being the most suitable reference state); an average of BOC & EOC insertions (30% in the example files) for cases used in the condensation of nuclear data. [Note that, in the example calculations, the cases producing nuclear data (& fluxes) in the 'perturbation' and 'shutdown margin' blocks all used rods fully withdrawn. This is correct for the 'perturbation' cases, since they are for the EOC state. The 'shutdown margin' cases are for BOC conditions, they should ideally be performed for an appropriate rod insertion - in practise, any difference should be negligible.]

Many (but not all) of the programs still use fixed format input: not only the position of numbers within each card, in some places specific numbers of blank cards are required to be input. (You have been warned!)

The example files of the following Section are all for the same core & fuel cycle calculation (there are exceptions, but they are not significant). Here is a brief description of the relevant details. A breederless reactor, with 2 core zones of different Pu enrichments. The detailed model represents a total of 52 different material zones, in 8 groupings: inner core - 8 axial zones, outer core - 3 radial x 8 axial zones for a total of 24 zones, 4 plenum zones, 4 plugged pin zones, 5 zones of rod follower or wrapper without pins, 2 zones of steel shield, 1 of B₄C shield, and 4 of rod absorber. Within each one of these 8 groupings the zones have the same number density (for a clean core); this provides 8 separate material zones in the 'iterative condensation' calculations, and the condensed nuclear datasets. The rod absorber is included as both heterogeneous and homogeneous cell models, giving a total of 9 entries in the condensed nuclear data. In the MOSES rod worth calculations, the above detailed model is used, but with the number of outer core zones reduced from 24 to 16. The fuel cycle is in 4 batches of 9.75 months.

6 The Computational Route - a Detailed Description.

For a full description of each input option, refer to the particular program's manual for further details. SLAROM and CITATION have manuals in English^(1,2), other programs have manuals only in Japanese (if at all). Appendices I to IV describe the input requirements of PENCIL, CITATION, PERKY and MOSES. The minor programs (e.g. JOINT) have their input described in the main text. Example JCL and data files are listed in Appendix VII.

A number of figures have been included, as an aid to understanding the calculational route. Figure 6.1 shows how the different blocks of the route fit together, and the main parameters calculated in each block. For the 6 blocks, each figure 6.2 to 6.7 shows the flow of data between calculation steps for one block; to aid clarity, some of the less significant datasets have been omitted; figure 6.5 includes the symbol conventions used in all 6 figures.

When using the route, the initial part of JCL and data file names - POC0B17 in the example files - will have to be altered to whatever FACOM uid is given to the user. You are of course free to completely change the filenames, they are only used here to help define where the same (or similar) files are used in different places. Note that other file names - principally program load modules, but also some standard data files (e.g. JENDL-3.2 70-group data) - represent fixed files used by the whole section, and they should not generally be altered.

6.1 Iterative Condensation block

6.1.1 PENCIL Calculation

The PENCIL calculation for the 'basic fuel cycle' is identical to this calculation.

A batch job run using JCL file [1], with Data files [1] & [2] as input (see appendix VII for example file listings).

File .OBHZHBC3.MIC18G3 is an 18-group 'micro' file produced by the S-J-C calculation in an 'iterative condensation'. A pre-existing data file is used to start the first iteration; in later iterations this file is then taken from the S-J-C run of the previous iteration.

Files .PQINT.DATA(APPLE) and .PQINT.DATA(TABLE) are temporary files transferred between the different program blocks. They can be used to provide summary table/graph plotting output.

File .PQINT.DATA(CIT) provides input to the next step, CITDENS. [For the 'basic fuel cycle' block, this file is also used as input to the 'shutdown margin' calculations.]

Data file [1] - .OPDEND.DATA(GDI1P) - is input to PENCIL which allows it to calculate core/breeder number densities from geometric data, and also provides parameters controlling the iterations of PENCIL. Appendix I gives details of the various parameters, various points of note are given here.

I have not run the program with breeder zones, so cannot give guidance on its effects.

For the 'iterative condensation', LAST=1 is necessary to ensure the same case in each iteration. For 'basic fuel cycle', a larger number is used, but not so high that CPU time can expire.

ICZ, OCZ (+ABZ, RBZ, IBZ) are the CITATION material zones for which number densities will be produced (see data file [2]).

PENCIL is set up to recognize 85 specific nuclides, in a particular order; the NUID number must be non-zero for any isotope actually present.

Lines beginning '&' are necessary block control commands; the lines beginning 'inner ..' & 'outer..' are comment lines.

The particular case shown has both ZrH and B₄C diluent replacing some fuel - it uses NZRH, VB4C and associated parameters.

Note that the F parameter should have been increased to values for 18 groups (this is not important, the 8'th to 18'th values should take the default values).

Data file [2] - .OPCITD.DATA(O975M165) - is the data for the CITATION run which is embedded within PENCIL. The data is in a series of blocks, identified by the lines reading '001', '002' etc. This is not quite the standard form of CITATION input: the number density data, block '020', excludes the core (& breeder) materials - zones 1 to 32 in this case - as they are calculated by the PENCIL routines. Whilst there is an English manual⁽²⁾ for CITATION, I include Appendix II describing some of the important features used.

Burn-up calculations use a detailed geometric model: 52 zones, 32 are core - see block '012'.

In this case, microscopic x-sections and number density data are input. The N.D. data uses nuclide identification numbers in the same format as PENCIL. Refuelling sequencing is set up for using a fixed feed fuel composition (i.e. no re-cycling). Although each core zone is represented as a smeared number density in flux calculations, within each zone the number densities of each refuelling batch are represented as a sub-zone for the burn-up calculations. 'Feed stream', 'makeup stream' and 'fuel movement' is defined once for each fuel zone; just one 'discharge stream' is defined (since no further use is made. I think it should be possible to use 1 'feed stream' and 1 'makeup stream' per enrichment zone.

[Note, the CITATION input also defines a zone class. I have maintained a simple correspondence of 1 zone to 1 zone class.]

6.1.2 CITDENS Calculation

The CITDENS calculation for the 'basic fuel cycle' is identical to this calculation.

A batch job run using JCL file [2].

The data file .PQINT.DATA(CIT) was created by the preceding PENCIL calculation.

The cross-section file - .OBHZHBC3.MIC18G3 - must be the same as used in the preceding PENCIL calculation.

The output file .@GDI1.DENS is the number density data for all timesteps during the burn-up calculation. It is read as input by RZOUT3, and by other programs.

6.1.3 RZOUT3 Calculation

An interactive procedure, run by a Macro.

The program extracts number density data from CITDENS output; it transforms the data to the SLAROM input format; it can combine material zones (calculating average number densities) to give the output for an altered set of zones.

To make the MACRO available within an interactive session, need to run the enablement file, JCL file [3] - .TSS.CLIST - by inputting the following commands at the command prompt:

```
STREAM COMMAND
EX TSS
LOGON
```

This executes the file named 'TSS' (of type 'CLIST') during the re-login which is initiated. You should only need this procedure again if either of JCL files [3] or [4] are changed.

The actual MACRO is run by typing 'RZOUT3' at the command prompt, this initiates the Macro, JCL file [4] - .TSSMAC.CLIST(RZOUT3). You have to follow this with certain typed input that the program requires.

It will be necessary to edit both JCL files [3] and [4] to handle the filenaming convention and uid that you employ.

The 2 typed input parameters that RZOUT3 requires, and prompts for, are:

```
'input' file identifier
'output' file identifier
```

if a null value is given for 'output', then it takes the same value as 'input'.

There are 3 files associated with running RZOUT3:
.@'input'.DENS the number density file created by CITDENS
.OPEDID.DATA('input'RZ) control data, data file [3]
.OPFORT.DATA('output'RZ) output file, data file [4]

6.1.4 Hand Edit

This step takes the SLAROM format number density data output by RZOUT3, data file [4], and produces the SLAROM data used in the S-J-C calculation, data file [5].

The editing procedure used to be carried out by hand but, especially following the introduction of a 2-region/2-temperature model for the fuel, it has become sensible to adopt an automatic editing program. A program (EDSLA) is being written that will be able to produce SLAROM data files (for either 1-region or 2-region fuel cell models), and so will replace this 'Hand Edit' step. A separate user guide and program description for EDSL A will be published⁽³⁾. For now I just describe the output file that has to be created.

Data file [4] - .OPFORT.DATA(GDI1RZ) - contains number density data in the format required in a 'SLAROM data' file; it was created by the foregoing RZOUT3 calculation. Each block of nuclide/number density data in SLAROM format is taken from the input file, removing the 'ZONE nn' line, and used as the basis for one material zone in the 'SLAROM data' file that is created. (Note that data file [4] is actually a 52 zone version for the 'basic fuel cycle' block, whilst the output data file [5] is for the 8 zone model of the 'iterative convergence'.)

Data file [5] - OPDEND.DATA(GDG0S) - is the 'SLAROM data' file that is created for use in the following S-J-C calculation. It consists of SLAROM input data for each material zone. Each fuel zone has the n.d. data adjusted to fit the 2-region representation; the rod absorber zone is modelled twice, the second time as a multi-region representation; all remaining zones are modelled as 1 (homogeneous) region, with n.d. data unaltered from the input values. For a full description of SLAROM data see the manual⁽¹⁾, the following gives a brief description of the important items:

PREP - codeword at start of new zone (for homogeneous zones the only codeword)

Title line

Control Data

field 1	no. of regions in cell; 1 if homogeneous, otherwise 2/6 for fuel/absorber
field 4	0/1 for homogeneous/heterogeneous case
field 5	use 0/1 - slab/cylindrical cell for homogeneous/heterogeneous
field 7	value >0 for JENDL-3.2 nuclear data
field 8	-ve value, SLAROM will include all micro. data in nuclear data PDS (0 - no micro. data included)

Control Data

field 1	temperature <u>or</u> temperature of non-fuel isotopes + 10 ⁴ {Manual description is incomplete}
---------	---

field 2 0 - homogeneous,
 1.3 - recommended for heterogeneous cell
 Temperatures fuel temperatures, 1 value per region: (for
 heterogeneous model with non-uniform
 temperatures, only present if
 'Control Data', field 1 > 10⁴)
 Nuclides no. of isotopes in each region
 Width thickness of each region
 n.d. ID and number density of each nuclide, data
 for each region starts on a new card
 PATH - codeword, only used for heterogeneous zones
 Title line
 Control Data
 field 2 no. of zones in cell, 2/6 - fuel/absorber
 field 5 manual says use same value as in 'PATH'
 block, (1); example fuel data I was given
 had value 0
 Geometry Data - the SLAROM manual refers to 'region',
 'zone' and 'material' model sub-divisions, in my
 calculations these were identical, all equivalent to the
 regions of the PREP data.
 field 1 3 sets geometry as cylindrical
 fields 2,3,5 the no. of regions, 2/6 - fuel/absorber
 fields 9,11,12,15 fuel values were taken from an
 example dataset; absorber calculation was
 made more detailed (it seemed prudent)
 Region Data mapping of 'material' data (input in PREP)
 to 'SLAROM regions' of the geometry [always
 1:1]
 Size radius of each cylindrical region of the
 cell model
 PIJF - codeword, only used for heterogeneous zones
 Control Data
 Convergence where input is zero, default value is used
 Data
 Control Datum 0 value stops activation cross-section
 input
 EDIT - codeword, only used for heterogeneous regions
 Control Data
 field 4 0/1 - all/sub-set of regions are included
 in averaging of cross-sections
 Sub-set Line only present if above item is '1'.
 1'st & last included regions (values
 repeat). Only used in rod absorber model,
 to exclude fuel surrounding absorber S/A in
 supercell.
 'NAME' - name given to zone by SLAROM and associated
 calculations (may be truncated to 5 characters).

The absorber heterogeneous cell model consists of 6
 cylindrical regions. The inner 5 regions comprise a single
 absorber S/A, with the 6'th region consisting of the
 homogeneous fuel composition, to a size corresponding to
 the fuel:absorber ratio in the core. (It is possible to
 run RZOUT3 again for every case, producing a single reactor
 average fuel composition. Rather than this extra

calculation, I have just taken the inner core n.d. from the equivalent of data file [4].) Region 5 is the S/A wrapper, steel of a volume taken from the rod follower zone. Region 4 is Na coolant, at a volume taken from the absorber S/A data. The remainder of the steel in the absorber S/A forms region 3, the sheath round the absorber. The absorber itself is region 1 - at its actual, not theoretical, density. Region 2 was the small remaining volume fraction - the gap between absorber and sheath: I filled this with Na (at greatly reduced density) to avoid any possible problems with an empty region. For the 5 inner regions, n.d. values were just those of the component material - Na, steel or B₄C, the latter reduced to its fraction of theoretical density.

The fuel heterogeneous cell model consists of 2 cylindrical regions. The inner region is just fuel material, with the radius of a fuel pellet. The outer region consists of all other materials in the fuel zone - steel wrapper and clad, Na coolant, diluent - of a radius such that the fuel pellet region has the same fractional volume as does the fuel in the S/A. The fuel pellet region has the same fraction of theoretical density as the actual fuel pellet, except that any central hole in the pellet is smeared into the fuel pellet region, reducing the density. The number density data from the RZOUT3 calculation has to be increased; for each cylindrical region a multiplier of the inverse of the zone volume fraction is used - thus maintaining the average density over the whole cell model.

The non-uniform temperature representation is used for the 2-region fuel cell model. I do not know exactly which are the fuel isotopes that the region dependant temperatures are applied to. There appear to be further difficulties with the non-uniform temperature model, I had problems trying to implement a 3-region - bore:fuel:other material - hollow pellet model.

To provide SLAROM data for an S-J-C calculation that produces 18-group nuclear data for shutdown temperatures, an extra Hand Edit is done, producing output that only differs from data file [5] in having a uniform 473 K temperature.

6.1.5 S-J-C Calculation

A batch job using JCL file [5], it is actually a series of linked calculations. A SLAROM cell calculation of macroscopic nuclear data; a CITATION flux snapshot calculation; JOINT condensations of nuclear data, and various interface routines (also part of JOINT).

The SLAROM step uses as input the data file [5] - .OPDEND.DATA(GDG0S) - created by and described under the foregoing Hand Edit.

File POC0H##.JFS3J3.Y9406 is the 70-group JENDL-3.2 data, input in a section standard data file (a fixed data file).

Macroscopic nuclear data, to be used by CITATION, are stored in the internal file designated &&XSG70 - this also includes the microscopic nuclear data which is to be condensed by JOINT; it is the size of this file that limits the number of material zones modelled in this calculation.

The CITATION step uses as input data file [6]. It also uses the macroscopic nuclear data of file &&XSG70 - this is in place of the microscopic nuclear data and number densities as used in the CITATION calculation within PENCIL.

Data file [6] - .OPCITD.DATA(PHA1N70) - is the geometric and other standard CITATION data. This is essentially the same as data file [2], but with certain differences.

As just a snapshot (i.e. no burn-up) calculation, data blocks '002', '018', '034', '036', '091' and '093' are not required (they are only for burn-up calculations). Blocks '004' and '005' represent a much reduced number of material zones.

Blocks '012' and '020' are not used since macroscopic nuclear data replaces the microscopic nuclear data and number densities of the PENCIL dataset.

A modified version of the CITATION input is used, allowing an easy interface with the JOINT routines and nuclear data from SLAROM. Data blocks '008' and '023' are non-standard input, as are the 2 lines before the title.

Initial Lines

Program name

A material name (from 'SLAROM data') - use 1'st (core) zone, it has all 85 (PENCIL) nuclides present.

Block '008'

line 2

field 1 no. of energy groups (minus sign indicates input from PDS file)

field 2 no. of energy groups for downscatter

field 3 no. of energy groups for upscatter

field 4 number of material zones

line 3+ ID of material used in block '005'

line 4+

field 1 name of material zone in SLAROM calculation (note, 5 characters)

line 5 (space)

line 6 material zone supplying fission spectrum data

Output from the CITATION step, to be used by JOINT, are placed in the internal file designated &&CIT.

There are 3 condensation and associated JOINT calculation steps. The extraction of fluxes from the CITATION output; the condensation of microscopic nuclear data (using the

extracted fluxes); the transformation of condensed nuclear data to a format used by PENCIL/CITATION. Data files [7], [8] and [9] provide control data used in the 3 steps.

The internal files &&XSG70 and &&CIT transfer output data from the SLAROM and CITATION steps of the calculation. A temporary file - .@GDG170.FLUX - holds the fluxes extracted from the bulk of CITATION output.

Condensed microscopic nuclear data are produced in 2 files: temporary file .@GDG118.PDSMIC , a 'partitioned dataset' (a library of various files), the format used by SLAROM; a single binary file .OBHZHBC7.MIC7GJ3 , the format used by PENCIL and CITATION. The 2 datasets also adopt the 2 different nuclide numbering schemes.

Data file [7] - .OPEDID.DATA(PFC1F) - lists those material zones for which fluxes are to be extracted

```

line 4
      field 1   number of energy groups
      field 2   number of zones/fluxes to extract
line 5+  names of material zones (note 5 characters)

```

Data file [8] - .OPEDID.DATA(PFC1C) - controls the actual condensation

```

line 4   number of groups after condensation
         {Note this particular file is for 7-group,
         not 18-group, condensation.}
line 5   lower energy group of each condensation
         band. {For 18 groups, standard data is -
         2, 4, 6, 8, 10, 13, 16, 19, 22, 25, 28, 31,
         34, 37, 40, 43, 46, 70.}

```

```

4 lines for each condensed nuclear data set
PDSIN    zone name of data to be condensed
PDSOUT   zone name of condensed data
USERPDS  zone name of flux used in condensation
Title line.

```

Note that the 2 absorber datasets (homogeneous and heterogeneous) from SLAROM are both condensed using flux from the same zone in CITATION.

Data file [9] - .OPEDID.DATA(PFC1M) - controls the transformation of condensed nuclear data between SLAROM (PDS) and CITATION (binary file) formats

The 1'st block inputs some data on all nuclides used in the PENCIL/CITATION data format:

```

line 2   name of 1'st fuel zone (1'st zone with all
         85 nuclides)
line 3   total number of nuclides
line 4+
      field 1   nuclide PENCIL ID number (in order)
      field 2   nuclide SLAROM ID number
      field 3   atomic weight
      field 4   energy per fission

```

Subsequent blocks - 1 for each material zone in condensed nuclear data:

```

line 1
      field 1   material zone name (note 5 characters)

```

field 2 no. of nuclides for which data is
 stored (all nuclides used for fuel
 zones)
 field 3 descriptive title
 line 2 PENCIL ID numbers of nuclides stored

To produce nuclear data condensed to both 7 and 18 groups, the whole S-J-C calculation can be repeated, with the condensation data of data file [8] differing as described above. Alternatively, the 3 final job steps - JOINT2, JOINT3 and MICRO - can be repeated at the end of the job: the step names will have to be changed, as will the names of temporary files created (to prevent duplication); also, in the first JOINT2 step, for file &&CIT alter '(OLD,DELETE)' to '(OLD,PASS)'. [Note, I have not actually used this second method, preferring to repeat the whole S-J-C calculation after I am satisfied Keff has converged.]

For the condensation of nuclear data in 7 groups at shutdown temperatures, the S-J-C calculation is repeated, but with the appropriately modified data replacing the SLAROM input of data file [5].

The CITATION format microscopic data file - .OBHZHBC7.MIC7GJ3 or its re-named equivalent - from the final iteration is used in various of the subsequent calculation blocks (e.g. 'basic fuel cycle').

6.2 Basic Fuel Cycle block

6.2.1 PENCIL Calculation

This is identical to the PENCIL calculation in the 'iterative condensation'.

A batch job run using JCL file [1], with the 18-group 'micro' file .OBHZHBC3.MIC18G3 being the condensed data produced by the final iteration of the 'iterative condensation'.

6.2.2 CITDENS Calculation

This is identical to the CITDENS calculation in the 'iterative condensation'.

A batch job runs using JCL file [2], with the 18-group 'micro' file .OBHZHBC3.MIC18G3 being the condensed data produced by the final iteration of the 'iterative condensation'.

The temporary data file used to control the calculation - .PQINT.DATA(CIT) - and the number density output file - .@GDI1.DENS - are used in some of the later calculation

blocks (e.g. shutdown margin).

6.2.3 MASSN Calculation

A batch job run using JCL file [6], this JCL file includes data.

File .GDI1.DENS is the number density multiple-entry dataset created by the above CITDENS calculation. (The change of name, removing '@', corresponded to copying the temporary file to permanent filespace.)

JCL file [6] - .OPJCL.CNTL(GDI1MASS) - contains the following data:

```

line 1      cycle length (days, integer); no. of cycles
            fuel is irradiated; pin clad outer diameter
            (mm); wire diameter (mm)
line 2+     total no. of zones in CITATION calculation;
            no. of nuclides modelled in each zone
line 3      no. of zones in inner core; zone ID no.'s
line 4+     no. of zones in outer core; zone ID no.'s
line 5      no. of zones in axial blanket ; zone ID
            no.'s [0 0 if no zones]
card 6      no. of zones in radial blanket; zone ID
            no.'s [0 0 if no zones]
line 7
            field 1-3 No. of S/As in inner core/outer core/
                    radial blanket
            field 4  No. of cycles stored on multiple n.d.
                    file [-ve if MOC included (i.e. cycle
                    length > 356 days)]
            field 5  Radial blanket irradiation (cycles)
            field 6  PENCIL ID of 238Pu isotope
            field 7  0/1 for half/full core height modelled
                    [note - example dataset has wrong
                    value, causes factor 2 error in all
                    masses.]
line 8,     title

```

The only output is the results (i.e. no file of data for use in further calculations).

(A modified version of the MASSN program, a more flexible and general version, is being produced; a separate user guide and program description⁽⁴⁾ are being produced.)

6.2.4 RZOUT3 Calculation

This is virtually identical to the RZOUT3 calculation of the 'iterative condensation' block, an interactive procedure run by the Macro RZOUT3.

The subsequent S-J-C calculation is to use the full detail (52 material zones) of the PENCIL calculation, so the input and output files will actually be of the forms shown in

data file [3] and data file [4], rather than with the slightly different data of the 'iterative condensation' calculations.

It is carried out twice, for both BOC and EOC steps of the final (6'th) iteration of the CITDENS burn-up calculation (ISTEP values of 16 and 17).

6.2.5 Hand Edit

This procedure is virtually the same as that carried out under the Hand Edit step of the 'iterative condensation' block. The input, from the above RZOUT3 calculation, has a larger number of material zones, hence the larger number of zones in the output, which is shown as data file [10] (with repetitive data deleted, to reduce the length of the printed file).

One change to the output file is that on the Control Data line of the PREP block, the 8'th field is altered to zero. (Since no condensation is to be done using this data, no microscopic data need be included in the output produced by SLAROM in the S-J-C calculation.)

A total of 5 different output data files are produced. The BOC and EOC versions of the input data from RZOUT3 give 2 standard cases. The EOC case also has 2 modified versions: with all Na (nuclide 11) n.d. set to 0.0 in fuel zones (Na void state); with fuel pellet regions in fuel zones having temperatures increased 500 degrees (for Doppler calculation). the BOC case has the modified version of a uniform 473 K temperature (shutdown temperatures).

(When it is available, the EDSL program⁽³⁾ can replace this 'Hand Edit' step.)

6.2.6 S-J-C Calculations

This is similar to the S-J-C calculation of the 'iterative condensation' block, but the JOINT steps of the calculation related to the nuclear data condensation are excluded.

A batch job run using JCL file [7], with Data files [10] & [11] as input.

File POC0H##.JFS3J3.Y9406 is the standard 70-group JENDL-3.2 nuclear data file (a fixed data file).

Macroscopic nuclear data, to be used by CITATION, are stored on the internal file designated &&XSG70. This does not include the microscopic nuclear data since no JOINT condensation is carried out, hence the much increased number of material zones is possible.

Data file [10] - .OPDEND.DATA(GDI1S) - was created by the above Hand Edit.

Data file [11] - .OPCITD.DATA(PEE1D700) - has no significant difference from data file [6], the equivalent data file in the 'iterative condensation' S-J-C calculation. The more detailed model alters the geometry data of blocks '004' and '005', and the material names of block '008' (and the pre-title data); some output options of block '001' line 3 are changed; a reactor power value is input, rather than taking the 1 MW default value.

Note that the JOINT program has to be re-compiled in program step DATAP2, to increase the array size so that it can handle the large model used in this calculation.

The calculations produce no further data files.

The S-J-C program is run once for each of the 4 or 5 datasets created by the previous Hand Edit.

6.3 Perturbation block

Unlike in the 2 previous blocks, a single SLAROM calculation is used to provide data used in a series of CITATION calculations, so they are submitted as individual jobs, rather than the combined S-J-C calculation.

The PERKY program version used differs from that of the (Japanese) manual⁽⁵⁾, in that it is modified to take data from the JOINT interface routines. Appendix III describes the data input required by PERKY, to the extent that I am familiar with.

6.3.1 RZOUT3 Calculation

This is virtually identical to the RZOUT3 calculation of the 'iterative condensation' block, an interactive procedure run by the Macro RZOUT3.

The subsequent SLAROM calculation uses a reduced number of material regions - 8 in each fuel enrichment zone and 6 for all other materials. To achieve this, the control input data differs from data file [3] in the following way:

```
NEWZN(1)=1,2,3,4,5,6,7,8,9,10,11,12,13,14,15,16,9,10,  
          11,12,13,14,15,16,9,10,11,12,13,14,15,16,  
          4*17,4*18,5*19,2*20,21,4*22,38*0,  
ZNGR(1)=16*1,4*2,2*3,  
ISTEP=17  
NZONE=52,22,
```

(The shield region immediately before the rod absorber includes B4C, so it is classed as an absorber region in RZOUT3. The ISTEP value is for EOC.)

The output differs in form from that of data file [4] only in the reduction of the number of blocks of data, from 52

to 22.

6.3.2 Hand Edit

This procedure is very similar to that carried out under the Hand Edit step of the 'iterative condensation' block (it will, similarly, be possible to replace this 'Hand Edit' step with the program EDSL, when it becomes available). The input, from the above RZOUT3 calculation, has a larger number of material zones than the equivalent 'iterative condensation' case of data file [4]. The 'SLAROM data' file that is created is shown in data file [12] (with repetitive data removed).

Data file [12] - .OPDEND.DATA(GDI3PRT) - is of the same basic form as its 'iterative condensation' equivalent, data file [5], but there are a number of significant differences:

The data is for use with the earlier versions of SLAROM and other programs, those in which 6-character region names are used. This required the number of nuclides per region to be limited to 34: this was achieved by removing those with a zero number density.

The data from RZOUT3, above, is for a different number of material regions than either previous equivalent calculation, so there are a different number of data blocks to the output. There is just one, heterogeneous, version of the rod absorber material region modelled.

After the 22 blocks of data corresponding to the data blocks of the input file, all 16 of the data blocks for the fuel regions are repeated, twice. These repeated blocks are modified: the first set, corresponding to the void state, has Na number densities set to zero; the second set, corresponding to the Doppler perturbation, has fuel pellet temperatures increased 500 degrees. In each case, the material region name is modified to be unique - these modified names are used in the later calculations of the perturbation block.

In the Control Data line of each PREP block, the 8'th field is altered to zero; except that in the first fuel region - including its 'void' and 'Doppler' versions - the -20 value is retained. Microscopic data, used in PERKY, are to be taken from this single region.

6.3.3 SLAROM Calculation

This is similar to the first part of the S-J-C calculation of the 'iterative condensation' block. However, different (6-character) program versions are used.

A batch job run using JCL file [8], with Data file [12] as input.
Data file [12] - .OPDEND.DATA(GDI3PRT) - was created by the above Hand Edit.

File POC0H##.JFS3J3.Y9406 is the standard 70-group JENDL-3.2 nuclear data file (a fixed data file).

Macroscopic nuclear data, to be used in following CITATION and PERKY calculations, are stored in a temporary file - .@GDI3P.PDS70G (replacing file &&XSG70). This only includes microscopic nuclear data for a single fuel region (including 'void' and 'Doppler' variants).

6.3.4 CITATION, 70 groups, 'normal'

This is similar to the 2'nd & 3'rd parts of the S-J-C calculation of the 'iterative condensation' block (the CITATION snapshot and JOINT condensation calculations), however, the final 'MICRO' step is omitted. Also, different (6-character) program versions are used, and the JOINT program had to be re-compiled to handle the large model of this calculation.

A batch job run using JCL file [9].

The CITATION step is basically the same as in the 'iterative condensation' block. It uses as input the 70-group macroscopic nuclear data file .@GDI3P.PDS70G created by the foregoing SLAROM calculation. Geometric type data is input in data file [13].

Data file [13] - .OPCITD.DATA(PIJ3P70N) - differs from the equivalent 'iterative convergence' block input file in the following ways:

Block '001' Line 2
Field 7 To output fluxes on FT34 (not as Manual)
Field 12 To calculate both direct & adjoint flux

Blocks '004' & '005' represent the more detailed model.

Block '008' has the increased number of material regions, with the change to 6-character region names (also in pre-title line).

Block '040' is data for the adjoint flux calculation.

Output from the CITATION step, to be used by JOINT, are placed in the internal file designated &&CIT.

The temporary flux file - .@GDI3P.FT34G70N - contains flux data used in later calculations in the 'perturbation' block, both direct and adjoint fluxes, in 70 groups, for the 'normal' core state.

The JOINT steps comprise just the extraction of fluxes from the CITATION output and the condensation of nuclear data to

18 groups. (Only PDS format data is required, so the 'MICRO' step is omitted.) They use as input the same effective cross-section data file - .@GDI3P.PDS70G - as the CITATION step, as well as the internal file &&CIT created by that step. The data files [14] & [15] are also used as input.

Data file [14] - .OPEDID.DATA(PIJ3PF) - is of the same form as data file [7] of the 'iterative convergence' block, defining those regions for which fluxes are extracted. The numbers and names of material regions are different, corresponding to those in the CITATION step. Note the 6-character region names.

Data file [15] - .OPEDID.DATA(PIJ3PC) - is of the same form as data file [8] of the 'iterative convergence', defining the condensation. The condensation is to 18 groups, the material regions are as in the CITATION step (the rod absorber material is only included the once). Note the 6-character region names.

A temporary file - .@GDI3P70.FLUXN - holds the fluxes extracted from the bulk of CITATION output, used for the condensation.

The temporary PDS file - .@GDI3P18.PDSMICN - contains condensed nuclear data used in later calculations in the 'perturbation' block, in 18 groups, for the 'normal' core state.

6.3.5 CITATION, 70 groups, 'void'

This is basically the same as the previous calculation, except that there is no condensation, and so no JOINT steps.

A batch job run using JCL file [10]. Geometric type data is input in data file [16], as is the 70-group macroscopic nuclear data file .@GDI3P.PDS70G created by the JOINT condensation of the foregoing CITATION calculation.

Data file [16] - .OPCITD.DATA(PIJ3PV70) - differs from the input of data file [13] in the following ways:

- Block '001' Line 2 Field 12 No adjoint calculation
- Block '008' has the fuel material region names altered to the 'void' variants (also in pre-title line).
- Blocks '023' & '040' are omitted.

The temporary flux file - .@GDI3P.FT34G70V - contains flux data used in later calculations in the 'perturbation' block, direct flux only, in 70 groups, for the 'void' core state.

(No &&CIT file is created.)

6.3.6 CITATION, 18 groups, 'normal'

This is basically the same as the previous 70-group, 'normal' CITATION calculation, except that there is no condensation, and so no JOINT steps.

A batch job run using JCL file [11]. Geometric type data is input in data file [17], as is the 18-group macroscopic nuclear data file .@GDI3P18.PDSMICN created by the JOINT condensation of the 70-group, 'normal', CITATION calculation.

Data file [17] - .OPCITD.DATA(PIJ3PADJ) - differs from the input of data file [13] of the first CITATION calculation of this block in the following ways:

Block '008' Line 2 No. of groups altered to 18
Block '023' - omitted.

The temporary flux file - .@GDI3P.FT34G18N - contains flux data used in later calculations in the 'perturbation' block, direct & adjoint fluxes, in 18 groups, for the 'normal' core state.
(No &&CIT file is created.)

6.3.7 CITATION, 70 groups, 'Doppler'

This is basically the same as the previous 70-group, 'normal' CITATION calculation.

A batch job run using JCL file [12].

The CITATION step is basically the same as in the previous 70-group 'normal' calculation. It uses as input the 70-group macroscopic nuclear data file .@GDI3P.PDS70G created by the foregoing SLAROM calculation. Geometric type data is input in data file [18].

Data file [18] - .OPCITD.DATA(PIJ3P70D) - differs from the input of data file [13] in the following ways:

Block '001' Line 2
Field 7 No FT34 flux output.
Field 12 No adjoint flux calculation.
Block '008' has the fuel material region names altered to the 'Doppler' variants (but not in pre-title line).
Block '040' - is omitted (no adjoint flux).

Output from the CITATION step, to be used by JOINT, are placed in the internal file designated &&CIT. No 'FT34' format flux file is created.

The JOINT steps are the same as in the previous 70-group, 'normal', calculation. They use as input the same effective cross-section data file - .@GDI3P.PDS70G - as the CITATION step, as well as the internal file &&CIT created by that step.

Two data files are input - .OPEDID.DATA(PIJ3PFD) and

.OPEDID.DATA(PIJ3PCD) - which differ from the equivalent files used in the 70-group 'normal' calculation, data files [14] & [15], only in replacing the fuel material region names by those for the 'Doppler' variant.

A temporary file - .@GDI3P70.FLUXD - holds the fluxes extracted from the bulk of CITATION output and used for the condensation.

The temporary PDS file - .@GDI3P18.PDSMICD - contains condensed nuclear data used in later calculations in the 'perturbation' block, in 18 groups, for the 'Doppler' core state.

6.3.8 PERKY, neutron lifetime

A perturbation-type calculation of prompt neutron lifetime, in 70 energy groups, using the PERKY program variant ABLE.

A batch job, run using JCL file [13]. Control data is input from data file [19], together with nuclear data and flux files created by the foregoing SLAROM and CITATION calculations.

Data file [19] - .OPEDID.DATA(PIJ3PNL) - provides the data for the PERKY neutron lifetime calculation. Nuclear data are input in just the macroscopic format, for each of the 22 material regions in the 'normal' CITATION calculations.

Other files used:

70-group 'normal' nuclear data: .@GDI3P.PDS70G

70-group 'normal' direct + adjoint fluxes: .@GDI3P.FT34G70N

6.3.9 PERKY, delayed neutron fraction

A perturbation-type calculation of delayed neutron fraction, in 18 energy groups, using the PERKY program.

A batch job, run using JCL file [14]. The PERKY load module is re-compiled, to allow a large model to be used. Control data is input from data file [20], together with nuclear data and flux files created by the foregoing SLAROM and CITATION calculations.

Data file [20] - .OPEDID.DATA(PIJ3PDNX) - provides the data for the PERKY neutron lifetime calculation.

Block '003' macroscopic nuclear data are input for each of the 22 material regions in the 'normal' CITATION calculations.

Block '002' microscopic nuclear data, all provided by the same material region, input for each of the 16 fuel regions, for the 8 nuclides producing delayed neutrons. Mapping of microscopic data to material regions has no data (i.e. blank) for non-fuel regions.

Block '005' delayed neutron fraction is taken as independent of energy of neutron causing fission.

These data are standard PNC values (Ref. 6)

Other files used:

18-group 'normal' nuclear data: .@GDI3P18.PDSMICN

18-group 'normal' direct + adjoint fluxes: .@GDI3P.FT34G18N

6.3.10 PERKY, Na void

An exact perturbation calculation of Na void, in 70 energy groups, using the PERKY program.

A batch job, run using JCL file [15]. Control data is input from data file [21], together with nuclear data and flux files created by the foregoing SLAROM and CITATION calculations.

Data file [21] - .OPEDID.DATA(PIJ3PEV) - provides the data for the PERKY Na void calculation.

Block '003' macroscopic nuclear data are input for 38 material regions: each of the 22 material regions in the 'normal' CITATION flux calculations, plus 16 fuel regions from the perturbed 'void' case.

Block '008' the mapping defines each 'normal' fuel region to be replaced by its 'void' equivalent.

Other files used:

70-group 'normal' nuclear data: .@GDI3P.PDS70G

70-group 'normal' direct + adjoint fluxes: .@GDI3P.FT34G70N

70-group 'Na void' direct flux: .@GDI3P.FT34G70V

Note that the PERKY load module used (Y9308) does not allow as large a calculation size as does that (Y9607) used in the following calculation.

6.3.11 PERKY, Doppler (by zone)

A 1'st order perturbation calculation of Doppler, in 18 energy groups, using the PERKY program.

A batch job, run using JCL file [16]. Control data is input from data file [22], together with nuclear data and flux files created by the foregoing SLAROM and CITATION calculations.

Data file [22] - .OPEDID.DATA(PIJ3PFD0) - provides the data for the PERKY Doppler (by zone) calculation.

As in the preceding calculation, except that the perturbed material region names are the 'Doppler' variants (rather than 'void'). (Minor data changes for no. of groups and 1'st order rather than exact perturbation.)

Other files used:

18-group 'normal' nuclear data: .@GDI3P18.PDSMICN

18-group 'Doppler' nuclear data: .@GDI3P18.PDSMICD

18-group 'normal' direct + adjoint fluxes: .@GDI3P.FT34G18N

6.3.12 PERKY, Doppler (by nuclide)

A 1st order perturbation calculation of Doppler, in 18 energy groups, with the contributions calculated for individual nuclides, using the PERKY program.

A batch job, run using JCL file [17]. Control data is input from data file [23], together with nuclear data and flux files created by the foregoing SLAROM and CITATION calculations.

Data file [23] - .OPEDID.DATA(GDI3PFD1) - provides the data for the PERKY Doppler (by nuclide) calculation.

Block '003' macroscopic nuclear data are input for 38 material regions: each of the 22 material regions in the CITATION flux calculations for the 'normal' state, plus 16 fuel regions from the perturbed 'Doppler' state.

Block '002' microscopic nuclear data, all provided by the same material region, input for each of the same 38 regions as the macroscopic data. Only 20 nuclides can have data input [presumably must include all perturbed nuclides and burnable nuclides, see block '008']. Map of microscopic data onto material regions includes all 38 regions.

Block '008' only 10 nuclides can be perturbed in any 1 run, need multiple runs to cover all nuclides [but several nuclides have no temperature dependence in JENDL-3.2, so can be excluded: ID's 946, 948, 950, 962, 963, 966, 967, 817, 857, 887, 897]. Unsure of 'burnable nuclide' category (guess it is those nuclides with changing number density), but limited to 10 nuclides. {Consider varying the nuclides chosen, as a sensitivity study.}

Other files used:

18-group 'normal' nuclear data: .@GDI3P18.PDSMICN

18-group 'Doppler' nuclear data: .@GDI3P18.PDSMICD

18-group 'normal' direct + adjoint fluxes: .@GDI3P.FT34G18N

6.4 Rod Worth block

6.4.1 PDSDUMP calculation

A short program that extracts data from PDS format nuclear data files, in this case fission spectrum data.

A batch job using JCL file [18].

The JCL file includes a source listing of the program,

which is compiled at run time; it also includes the control data. The calculation uses as input a PDS nuclear data file, the 18-group file created in the final cycle of the 'iterative convergence' block - .@GDG118.PDSMIC . JCL file [18] is shown using a different PDS data file, taken from a calculation for the 'shutdown margin' block.

In practise it is only necessary to do one of the PDS DUMP calculations, from this and the 'shutdown margin' blocks, since there will be very little or no difference in fission spectra.

[To condense an 18-group fission spectrum to 7 groups, simply add values where groups are combined.]

6.4.2 RZOUT3 calculation

This is virtually identical to the RZOUT3 calculation of the 'iterative condensation' block, an interactive procedure run by the Macro RZOUT3.

The subsequent MOSES calculation uses a reduced number of material regions - 16 rather than 24 in the outer fuel enrichment zone and 6 for all non-fuel materials. To achieve this, the control input data differs from data file [3] in the following way:

```
NEWZN(1)=1,2,3,4,5,6,7,8,9,10,11,12,13,14,15,16,17,18,  
19,20,21,22,23,24,17,18,19,20,21,22,23,24,4*25,4*26,  
5*27,2*28,29,4*30,38*0,  
ZNGR(1)=24*1,4*2,2*3,  
ISTEP=16  
NZONE=52,30,
```

The output differs in form from that of data file [4] only in the reduction of the number of blocks of data, from 52 to 30.

The calculation is carried out for just the BOC step (ISTEP=16).

6.4.3 Hand Edit

This step takes the SLAROM format number density data output by the above RZOUT3, data file [4], and produces the data used in the following MOSES calculations, as shown in data file [24] (with repetitive data removed).

The editing procedure has been carried out by hand, but it would be advantageous to have a program to automatically alter the number density data to the format used by MOSES. (Such a program, EDMOS, is being written, a separate user manual and program description⁽⁷⁾ will be published.) For now I just describe the output file that has to be created by hand. Appendix IV describes the format of MOSES input.

Data file [24] - OPMOSD.DATA(GDI1DY) - is the 'MOSES data' file used in the following calculation.

There are actually 4 different datasets created, with small differences: AXIAL & ASYGROUP data include absorber rod S/A types for 0%, 50% and 100% rod insertions; the cases differ in the S/A types assigned to rod positions in the REFMAP data.

LOADING 1 invokes the 1'st set of S/A loadings in REFMAP (the whole core) - in these steady state calculations, this is the only LOADING used.

360° symmetry is necessary for (some) cases with 1 rod stuck, it is easiest to retain this geometry in all cases.

I inadvertently used NAZN>14 when setting up the model - no problems were encountered.

Uses hex-Z geometry with interpolation of pseudo-tri-Z ratings.

DELDAY value input even for steady state calculation. In REGION, rod absorber material is given homogeneous version of nuclear data.

The fission spectrum data in MICROXS come from the previous PDSDUMP calculation.

XSNUC lists only those PENCIL nuclides actually present in the model.

Minor Actinides 21 to 28 are set to nuclide class 6, not 3 (following an example dataset I was given).

REACT1 & YIELD are supposedly necessary for steady state calculations (I would have thought that the presence of capture and/or fission cross-section values in the nuclear data file would be sufficient to ensure that these reactions are included in the Keff calculation).

The ATDENS number density data contains, for each material region, a value for each nuclide included in the MOSES calculation (0 if not present) - this is obtained by editing the RZOUT3 output file.

Even though this is a steady state calculation, a refuelling period must be given with the REFUEL data. REFMAP need not assign a S/A type to each S/A in all the NLAY rings, the radial boundary conditions will be applied at the edge of those S/As assigned by REFMAP. For the 4 different cases, just the assignments to S/A types 5, 6 and 7 require re-arranging.

6.4.4 MOSES calculation

Steady state calculations for different rod insertions; Keff differences are used to calculate nominal values of rod group worths.

A batch job using JCL file [19].

Input is the 'MOSES data' file [24] created in the above Hand Edit.

File .OBHZHBC7.MIC7GJ3 is the 7-group microscopic nuclear data file created in the 'iterative convergence' block.

If the effect of shutdown temperatures on nuclear data is being calculated using MOSES (rather than S-J-C), then one of the above MOSES calculations is repeated, but with the nuclear data file replaced with that created for cold shutdown conditions (from the 'iterative convergence' block).

6.5 Shutdown Margin block

There are 2 separate groups of calculations, calculating the reactivity effects of increasing the core size, and of increasing material densities.

Appendix V shows how the results of the calculations are combined to produce the core expansion reactivity change; this result combines with the reactivity effect of temperature on nuclear data (from previous MOSES or S-J-C calculations). Appendix VI describes the combination of this and other factors to produce values of shutdown margin.

The following is the route for calculating the reactivity effect of increasing core size.

6.5.1 Hand Edit

This procedure takes a 'CITATION data' file, data file [11], and edits it to create the 'CITATION data' file used in the following S-J-C calculation.

Data file [11] - .OPCITD.DATA(PEE1D700) - is the 'CITATION data' file that was used in the 'basic fuel cycle' block.

The output file differs only slightly from the input file - Within data block '004', either the axial or the radial mesh intervals are all increased in size by a set small fraction (I used a factor 1.01). Two datasets are created, one each for axial and radial mesh adjustments.

6.5.2 S-J-C calculation

A batch job virtually identical to the S-J-C calculation of the 'basic fuel cycle' step. The JCL differs from that of JCL file [7] only in having the 'CITATION data' file .OPCITD.DATA(PEE1D700) replaced by the data file created in the above Hand Edit.

The calculation is performed for both of the data files created by the above Hand Edit.

The material density reactivity coefficient calculations

start here. The calculations are rather similar to the 'perturbation' block, except that the model is simplified to a single material region for each Pu enrichment zone, and the '5-character' versions of the programs are used.

6.5.3 RZOUT3 calculation

This is completely identical to the RZOUT3 calculation in the 'iterative condensation' block: if a copy of the output file from that calculation is available, then the step does not need repeating here.

If the calculation has to be repeated, it uses the same input files as previously: the number density file - .@GDI1.DENS - created by the 'iterative condensation' CITDENS calculation, and 'RZOUT3 data' file that produces the least detailed model (NZONE=52,8,). The output that is produced will, again, be of the same format as data file [4], but for a lesser number of material regions.

6.5.4 Hand Edit

This step takes the SLAROM format number density data output by RZOUT3, similar to data file [4], and produces the SLAROM data used in the S-J-C calculation, data file [25].

The editing procedure is simple enough to make any automated process superfluous.

Data file [4] - .OPFORT.DATA(GDI1RZ) - contains number density data in the format required in a 'SLAROM data' file; it was created by the foregoing RZOUT3 calculation. (Note that data file [4] is a 52 zone version for the 'basic fuel cycle' block; the 'shutdown margin' block actually uses a version with an 8 zone model.)

Data file [25] - OPDEND.DATA(GDI1RZEC) - is the 'SLAROM data' file that is created for use in the following calculation. It consists of SLAROM input data for each material zone.

Each zone is modelled as a 1-region homogeneous model (the rod absorber is only represented the once).

Microscopic data is included for all isotopes of all material regions.

All 8 material regions are repeated, as Doppler versions - the material region names are altered, only fuel region temperatures are changed.

(When it is available, the EDSL A program⁽³⁾ can replace this 'Hand Edit' step.)

6.5.5 SLAROM calculation

This is similar to the first part of the S-J-C calculation of the 'iterative condensation' block.

A batch job run using JCL file [20], with Data file [25] as input. (The JCL differs from that for the 'perturbation' block in using a 5-character program version.)

Data file [25] - .OPDEND.DATA(GDI1RZEC) - was created by the above Hand Edit.

File POC0H##.JFS3J3.Y9406 is the standard 70-group JENDL-3.2 nuclear data file (a fixed data file).

Macroscopic nuclear data, to be used in following CITATION and PERKY calculations, are stored in a temporary file - .@GDI1P.PDS70G (replacing file &&XSG70). This includes both 'normal' and 'Doppler' state variants; microscopic nuclear data are included for all material regions.

6.5.6 CITATION, 'normal'

This is similar to the 2'nd & 3'rd parts of the S-J-C calculation of the 'iterative condensation' block, the CITATION snapshot and JOINT condensation calculations. (The JCL differs from that for the 'perturbation' block in using 5-character program versions.)

A batch job run using JCL file [21].

The CITATION step is basically the same as in the 'iterative condensation' block. It uses as input the 70-group macroscopic nuclear data file .@GDI1P.PDS70G created by the foregoing SLAROM calculation. Geometric type data is input in data file [26].

Data file [26] - .OPCITD.DATA(PFC3N70X) - differs from the equivalent 'iterative convergence' input data file in only one particular: the main rod absorbers are modelled fully withdrawn (rather than at the average insertion used in the 'iterative convergence' calculation), thus one less axial material zone is modelled in data blocks '004' & '005'.

Output from the CITATION step, to be used by JOINT, are placed in the internal file designated &&CIT.

There are 3 condensation and associated JOINT calculation steps. The extraction of fluxes from the CITATION output; the condensation of microscopic nuclear data (using the extracted fluxes); the transformation of condensed nuclear data to the CITATION format. Data files [27], [28] and [29] provide control data used in the 3 steps.

The same effective cross-section data file as in the CITATION step - .@GDI1P.PDS70G - is used as input, as well as the internal file &&CIT created by that step. A

temporary file - .@GDI1P70.FLUXN - holds the fluxes extracted from the bulk of CITATION output.

Condensed microscopic nuclear data are produced in 2 temporary files: file .@GDI1P18.PDSMICN, in the PDS format used by SLAROM; file .@GDI1P18.MICRON in the binary format used by CITATION. The 2 datasets use the 2 different nuclide numbering schemes.

Data file [27] - .OPEDID.DATA(PFC3FX) :
Data file [28] - .OPEDID.DATA(PFC3C18X) :
Data file [29] - .OPEDID.DATA(PFC3MX) . These differ only slightly from the equivalent data files - [7], [8] and [9] - from the 'iterative condensation' block. There is only a single rod absorber material, region name 'AHOMO'. The condensation is to 18, not 7, groups.

6.5.7 Hand Edit

Data files used in the above CITATION calculation are edited to provide data for the following CITATION. The only changes made are to the material region names - exchanging them for the equivalent 'Doppler' variants from the foregoing SLAROM calculation. The files affected are as follows:

.OPCITD.DATA(PFC3N70X)	→	(PFC3N70D)
.OPEDID.DATA(PFC3FX)	→	(PFC3FD)
.OPEDID.DATA(PFC3C18X)	→	(PFC3C18D)

6.5.8 CITATION, 'Doppler'

This is almost the same as the previous CITATION calculation. The main difference is that the data used is for the 'Doppler' state; also, the 'MICRO' step of the JOINT part of the calculation is omitted.

A batch job, run by JCL file [22].

The calculation uses as input the 70-group macroscopic nuclear data file .@GDI1P.PDS70G created by the earlier SLAROM calculation. Also used are the 3 data files created by the foregoing Hand Edit: geometric 'CITATION data' in file .OPCITD.DATA(PFC3N70D); flux extraction data in file .OPEDID.DATA(PFC3FD); condensation data in file .OPEDID.DATA(PFC3C18D).

Output from the CITATION step, to be used by JOINT, is placed in the internal file designated &&CIT. Temporary file .@GDI1P70.FLUXD holds the fluxes extracted from the bulk of CITATION output. The condensed nuclear data, in SLAROM (PDS) format, is output in temporary file .@GDI1P18.PDSMICD.

6.5.9 PDSDUMP calculation

A short program that extracts data from PDS format nuclear data files, in this case fission spectrum data. This is the same as the PDSDUMP calculation of the 'Rod Worth' block.

A batch job using JCL file [18].

Data is input as a PDS nuclear data file, that created in the foregoing 'normal' CITATION calculation -
.@GDI1P18.PDSMICN .

6.5.10 PERKY Package calculation

This was provided to me as a pre-assembled JCL script, which linked a CITATION calculation of direct & adjoint fluxes with PERKY calculations of density perturbations for fuel, coolant and structural core materials. A pre-processor links the various data files together, and a post-processor calculates values of material density reactivity coefficients.

A batch job using JCL file [23]; the JCL includes one data line with 3 values that specify a n.d. dataset to be used -
field 1 No. of cycle used (final cycle in CITDENS)
field 2 Step within cycle: 1/2/3 - BOC/MOC/EOC
(cycle > 1 year); 1/2 - BOC/EOC (cycle < 1 year)
field 3 (same as field 2)

Various 18-group nuclear data files created in the foregoing CITATION calculations are used -
.@GDI1P18.PDSMICN PDS format, 'normal' state
.@GDI1P18.PDSMICD PDS format, 'Doppler' state
.@GDI1P18.MICRON binary format, 'normal' state.

File .@GDI1.DENS is the number density dataset created by the CITDENS calculation of the 'basic fuel cycle' block; file .@GDI1.CITDAT is the temporary data file used in the same CITDENS calculation [file renamed from .PQINT.DATA(CIT)].

A whole series of further input data files are required. There are data files [30], [31], [32] and [33]. Delayed neutron data, in 18 groups, is input in a binary file - .DELAY.DATA - which is a copy of a standard section data file (a fixed data file). Three files of post-processor control data are input - .PQINT.DATA(POFUL) , (POSTR) , (POCOL) - consisting of:
line 1 value 2/3/4 indicating fuel/structure/coolant
line 2 value 1

Data file [30] - .OPEDID.DATA(NAMEF#XX) - gives the name of each material zone in order, first for the 'normal' state, then those for the 'Doppler' state.

Data file [31] - .OPEDID.DATA(IDNUMXX) - gives the PENCIL ID no. then the SLAROM ID no. of all nuclides in the nuclear data sets.

Data file [32] - .OPEDID.DATA(CHIMG) - gives the 18-group fission spectrum data from the previous PDS DUMP calculation.

For each fissioning nuclide (i.e all actinides with non-zero n.d.) -
 nuclide ID no. (SLAROM version)
 fission spectrum (18 values)
End with value 0.

Data file [33] - .OPEDID.DATA(ZONE#XX) - defines the zones for which the perturbation results are output.

Line 1 title line (do not alter)
Line 2+
 field 1 CITATION ID no. of 1'st of a
 consecutive set of 'volumes'
 field 2 CITATION ID no. of last of a
 consecutive set of 'volumes'
 field 3 title given to this zone (used in
 output)

Note that, as originally supplied, the PERKY package also included calculations of Doppler, Na void, beta-effective and neutron lifetime. I deleted the sections that carried out these calculations, the 3 reactivity coefficients being all I required from this calculation (note that this calculation would be inadequate for calculating Na void, because of the use of only 18 groups).

6.6 3D Burn-up block

6.6.1 S-J-C Calculations

It is up to you whether you use S-J-C as a quick route to find approximate values for BOC and EOC rod insertions that give balanced ratings and the required Keff value. The calculation will be virtually identical to the S-J-C calculation in the 'basic fuel cycle' block, with JCL taken from the JCL file [7]. The only difference is that the 'CITATION data' file .OPCITD.DATA(PEE1D700) (data file [11]) has to be changed. The data file only requires a minor modification, altering block '005' to change the rod insertions. (It may also be necessary to modify blocks '004' and '005' to insert extra axial layers, to allow rod insertions to flux meshes that don't correspond to material zone boundaries.)

6.6.2 Hand Edit

This produces data file [34] - .OPMOSD.DATA(GDI1MBRN) - the input data file for the following 3D burn-up MOSES

calculation. It is basically the equivalent data file from the 'rod worth' MOSES calculation - data file [24] - but with the following changes.

A burn-up, rather than a steady state, calculation.

The same as the PENCIL calculations, models repeated irradiation/ refuelling cycles until it approximates to fuel cycle equilibrium. Individual S/A refuelling batches are modelled explicitly.

Use the highest level of symmetry possible, to reduce calculation time - in this case 120° symmetry.

Note some changes to iteration control parameters NITR, NIIT and EPSF, to values suited to transients.

Each irradiation step is a half cycle, DELDAY - BOC to MOC, then MOC to EOC.

In AXIAL and ASYGROUP, each rod insertion to be used defines a separate S/A type; each fuel S/A type is defined twice - the 1'st time is used for the original loading, the 2'nd time for refuellings.

The axial mesh used for irradiated fuel is defined in AXIAL (all fuel S/As are originally of axially uniform composition).

The rod absorber material uses heterogeneous nuclear data.

In REACT1 & YIELD, the limitation on the number of entries meant that the burn-up chain could not include the full detail of minor actinides included in the PENCIL/CITATION calculations. A minimal series was adopted; fission product and ¹⁰B reactions were included (the latter only in diluent in the fuel, not in absorber rods).

The number density data is for the clean core - hence only 1 dataset per core Pu enrichment zone. The values can be taken from the temporary CITDENS data file - .PQINT.DATA(CIT) - from the 'basic fuel cycle' block (the same file is used in the 'shutdown margin' block, renamed as .@GDI1.CITDAT).

The REFUEL & REFMAP data define 6 sets of S/A loadings. First is the initial core loading (note the different S/A types c.f. later loadings). The 2'nd loading alters just the main rods, setting them to EOC insertions. The remaining 4 loadings each represent the refuelling of 1 of the 4 explicit fuel batches; rods are also moved to BOC positions.

After the PLANT & CRPOS data, a whole series of further calculation steps are defined; the calculation has a total of 9 cycles to reach equilibrium and then allow one cycle for each batch. There are 3 steps per cycle:

- (1) refuel/initial loading (rods to BOC positions)
 - flux calculation (BOC)
 - burn-up calculation (BOC to MOC).
- (2) move rods to EOC position [LOADING 2]
 - flux calculation (MOC)
 - burn-up calculation (MOC to EOC).
- (3) (no loading) [LOADING 0]
 - flux calculation (EOC)
 - (no burn-up) [DELDAY = 0.0].

One set of REUEL & REFMAP data can be used repeatedly, by re-using the appropriate ID number in LOADING. Note that parameters are always as in the previous calculation step, unless overwritten.

6.6.3 MOSES calculation

A transient consisting of repeated burn-up and refuelling until fuel cycle equilibrium is approach (the same as the earlier PENCIL calculations). This pseudo-tri-Z model provides a reasonably accurate assessment of peak pin ratings.

A batch job using JCL file [24]. Compared with the 'rod worth' MOSES calculations, the JCL parameters N10, NX, NS are increased to accommodate the larger files needed by the transient calculation - all file size parameters of this type depend on the model size.

Input is the 'MOSES data' file [34] created in the above Hand Edit.

File .OBHZHBC7.MIC7GJ3 is the 7-group microscopic nuclear data file created in the 'iterative convergence' block.

The calculation is repeated as necessary, to adjust the BOC & EOC rod insertions to give both flat radial rating profiles and acceptable values of Keff (the latter is none too important, because of uncertainties in Keff associated with 7-group nuclear data).

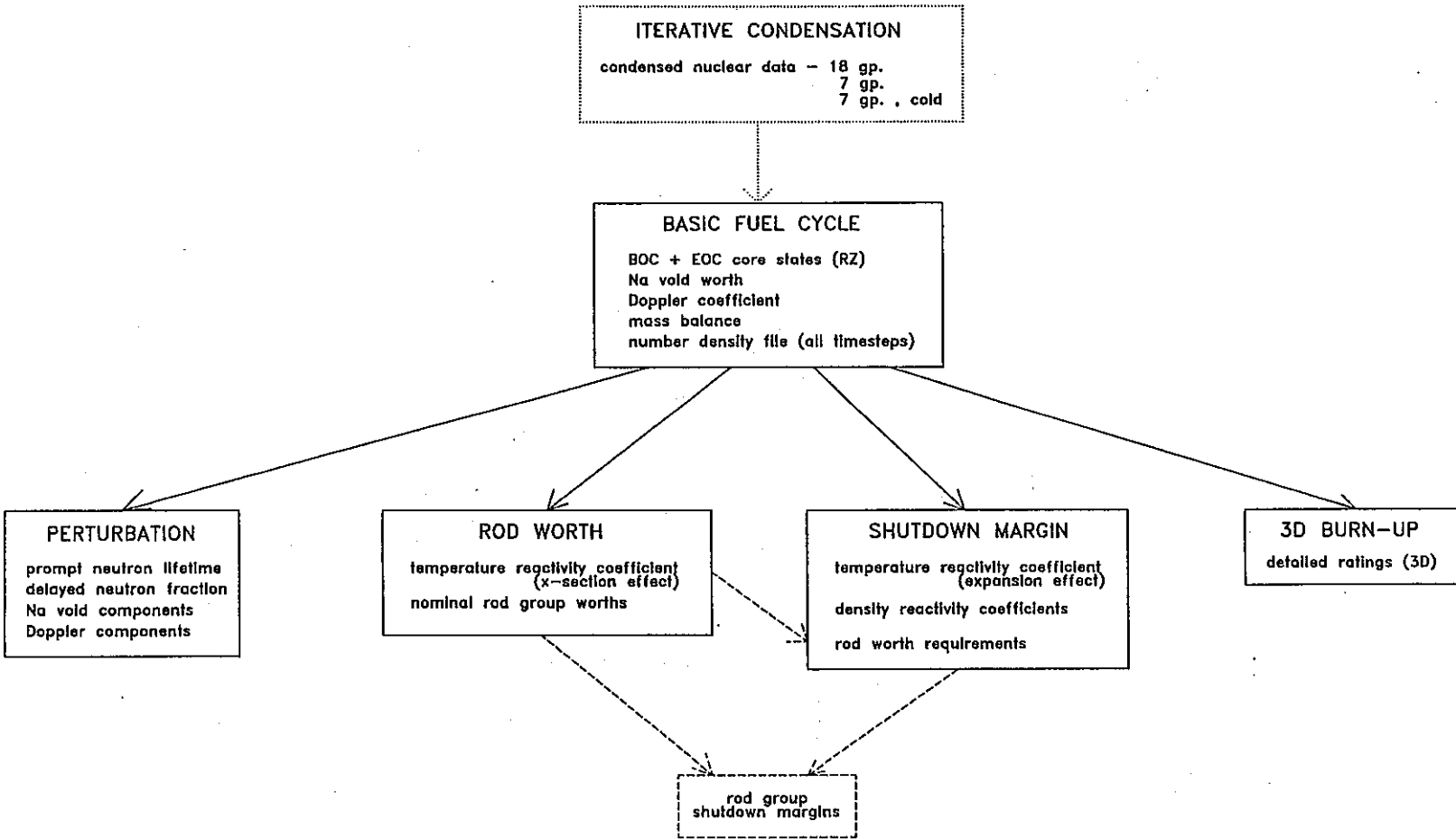


Figure 6.1 Calculation route, with parameters and major data items created by the separate blocks

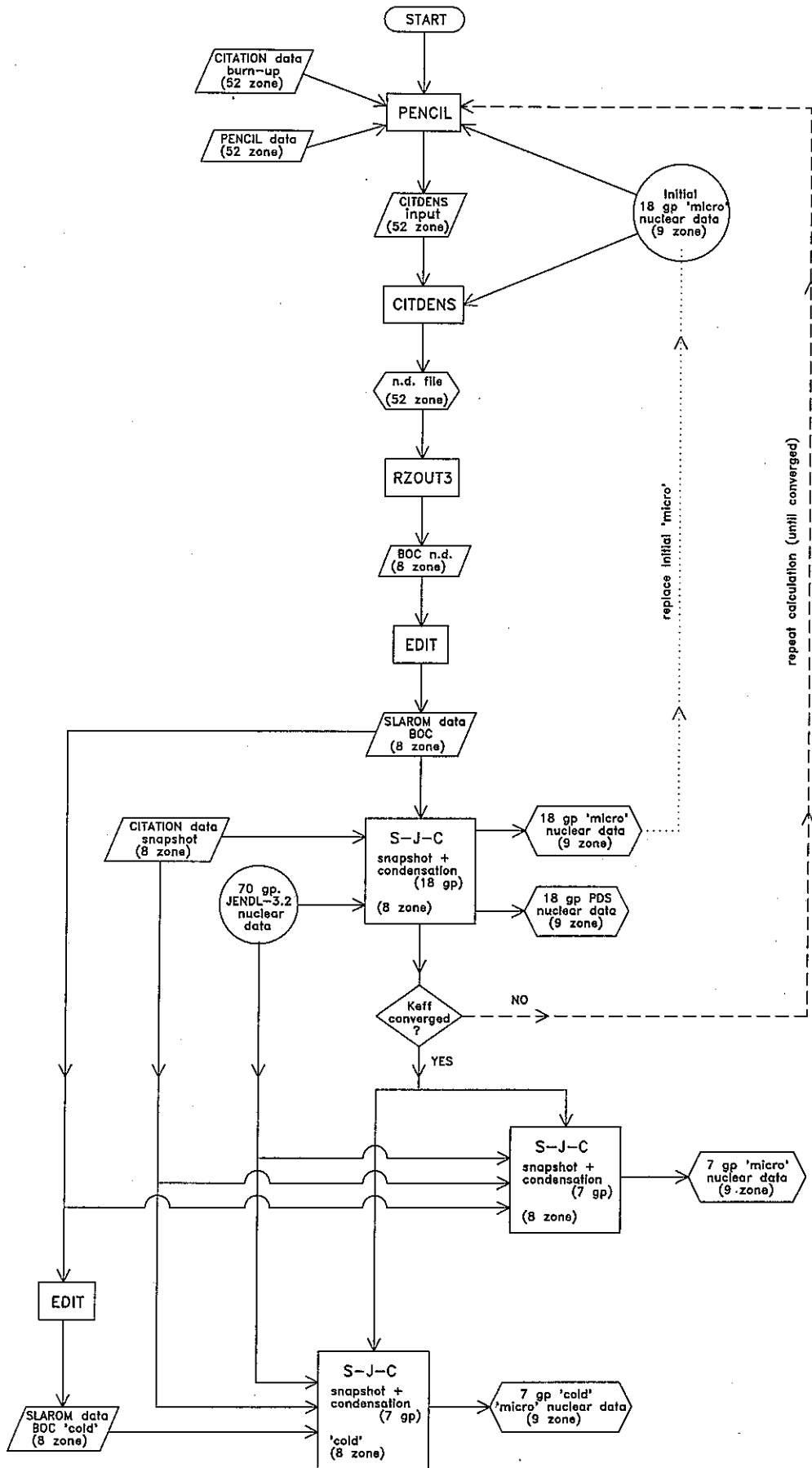


Figure 6.2 Schematic diagram - Iterative Convergence

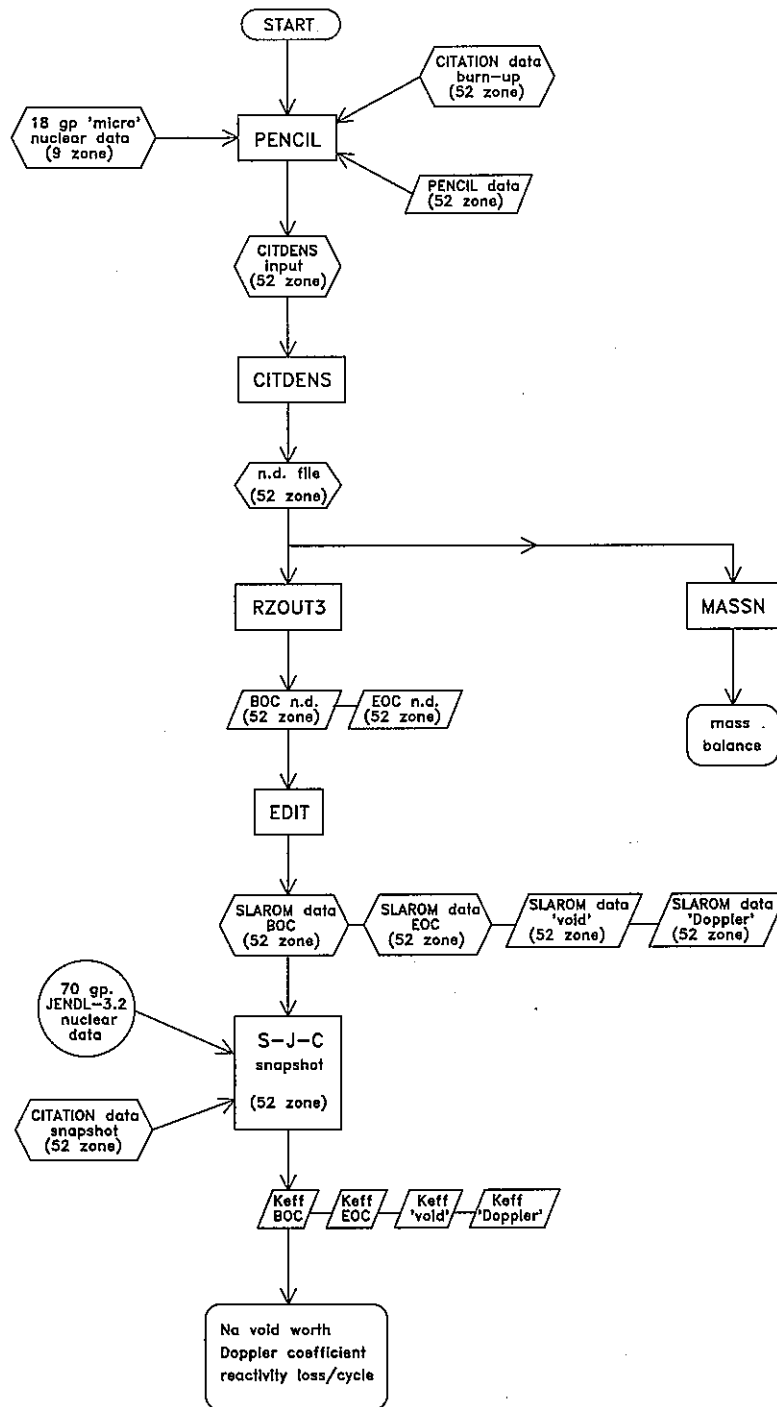


Figure 6.3 Schematic diagram - Basic Fuel Cycle

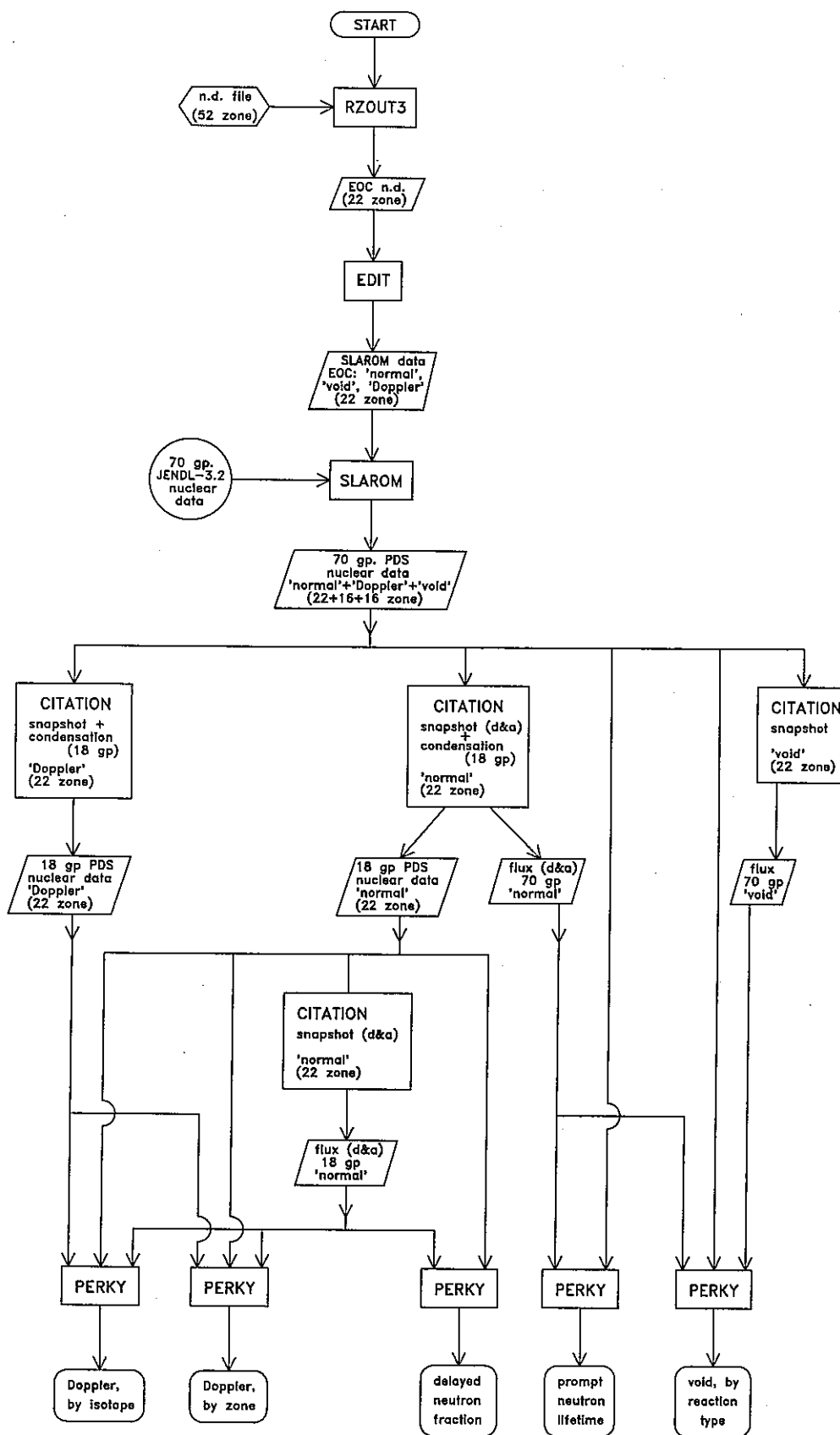


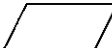
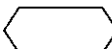




Figure 6.4 Schematic diagram - Perturbation

SYMBOL CONVENTION (figs. 6.2 to 6.7)

-  program, or similar calculation step
-  start point, or final result of calculation
-  data created just for/by the current block
-  data transferred from or required in another calculation block
-  fixed data file
-  several similar datafiles, from/for repeated runs of preceding/following steps

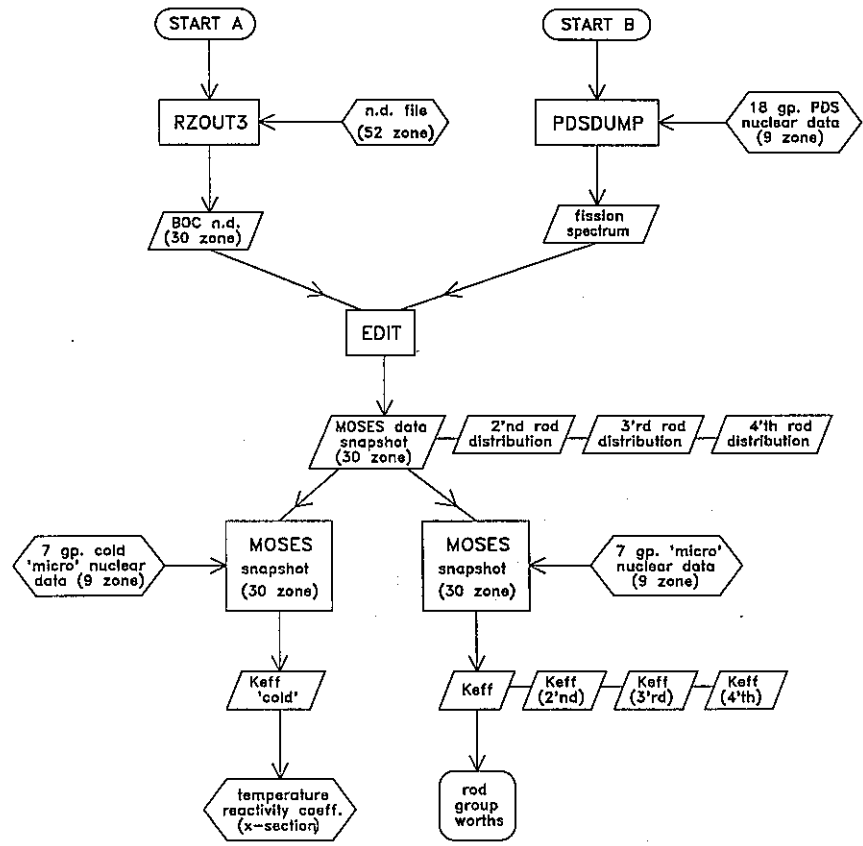


Figure 6.5 Schematic diagram - Rod Worth

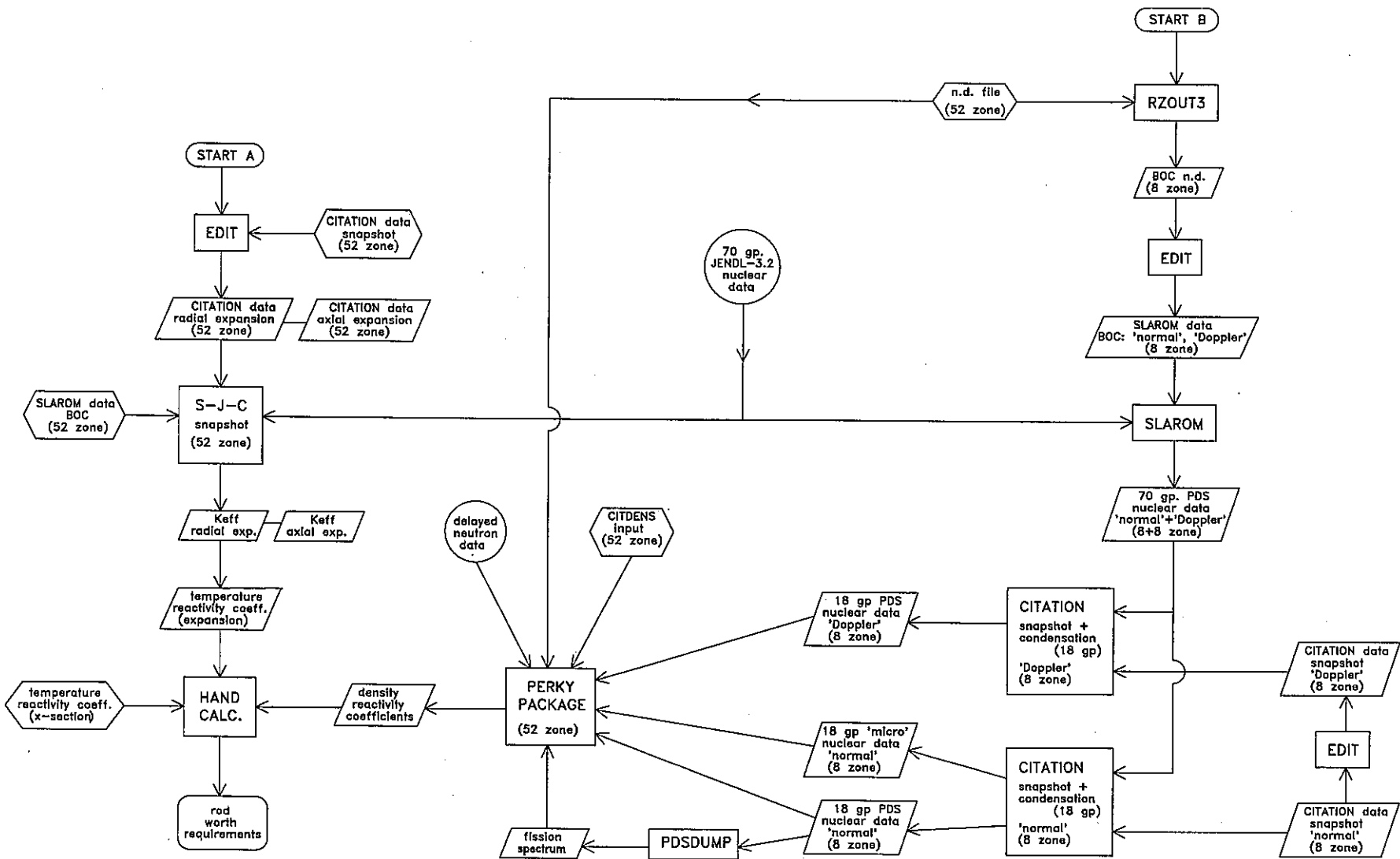


Figure 6.6 Schematic diagram - Shutdown margin

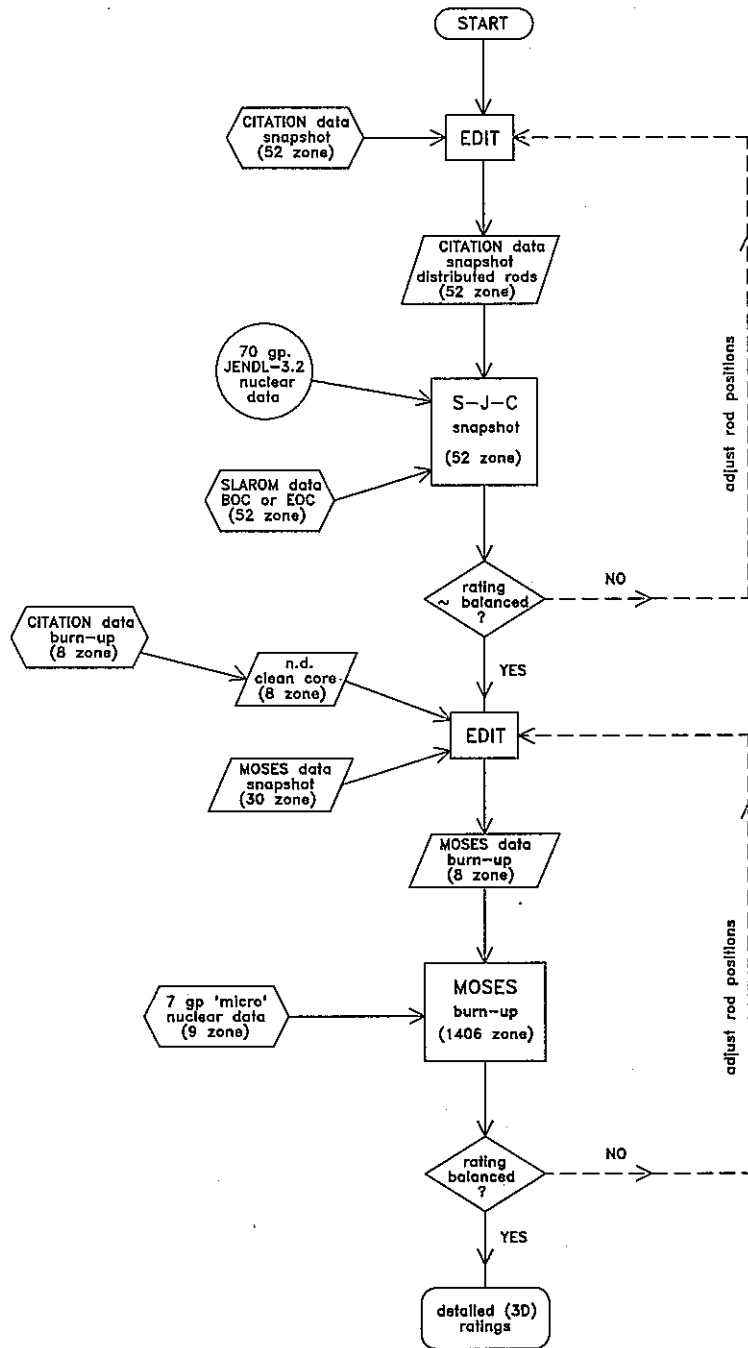


Figure 6.7 Schematic diagram - 3D burn-up

7 The Computational Route - Future Modifications.

Concurrent with the production of this description of the calculational route, a number of computer programs are being written to update the route. These programs will have no impact on the structure of the calculational route, simply replacing steps in the route as described above.

Two programs, EDSL^A(³) and EDMOS(⁷), will be available to replace the Hand Edit steps that take the number density data extracted by the RZOUT3 program and transform it to the formats required by SLAROM and MOSES, respectively. These programs will be advantageous in removing processes that are time-consuming (and potentially error-prone), especially when the 2-region fuel model is used in SLAROM. User guides are being produced for these programs(^{3,7}).

A modified version of the MASSN program is being produced. The main change is that it will have the flexibility to handle cores with more than 2 core and breeder zones (such as BN600) with varying numbers of irradiation batches. A user guide is being produced(⁴). [To allow the full calculational route to handle such cores, PENCIL will also have to be modified - this has yet to be attempted.]

Appendix I - PENCIL input data

VARIABLE	VALUE	DESCRIPTION
LAST	(max 30)	No. of Pu enrichment iterations. +ve/-ve - power plot output is core+axial blanket/core; 0 - just calculates N.D.s , no CITATION
PURICH (i≤2)		inner/outer Pu enrichments, (w/o) (initial guess, if LAST>1)
LPOWER	[430]	pin rating limit, W/cm (0.- not used)
TPOWER	[2600]	reactor power, MWth , assumes fixed thermal efficiency (EPOWER*26) (only 1 of this and EPOWER needed)
EPOWER	[100]	reactor power, 10's MWe , ignored if TPOWER used (TPOWER/26)
DAY(i≤3)	[365;3;4]	cycle length (days); No. of batches for fuel; No. of batches for radial blanket [max. of last 2 is basis for no. of cycles calculated, overriding value in CITATION input file]
COREH (i≤6)		axial height (cm) of each core zone (from top to bottom) [I used 8 axial zones, without problem]
ICZ(i≤24)		zone no's of inner core
OCZ(i≤24)		zone no's of outer core
ABZ(i≤24)		zone no's of axial blanket
RBZ(i≤24)		zone no's of radial blanket
IBZ(i≤24)		zone no's of internal blanket, (can be any position)
NUID (i≤99)		map of isotopes present in core (0 - absent; i - present)
F(i≤20)	[1;1; .94734;4*0] 7 gps [7*1; .8068;10*0] 18 gps	map of which energy groups are used in output flux values (usually limited to 100 kev) (does not affect calculation)
BOND	[0]	can filled with 0/1 - He/Na
DUCTLS (i≤3)	[0;6; 1.19277]	i=1, 0/1 with/w'out S/A wrapper i=2, no. of tie-rod pins } only if i=3, pin pitch/pin diameter} i=1 is 1
RONA	[.849052 @ 430 C]	Na density (g/cc)
ICNROD	[0]	0 - rods always out; 1 - split cycle in 2, rods 50%/0% in, needs full height model (see CITATION for rod movement)
DFUEL		fuel pellet outer diameter (mm) [only 1 of this & SMEAR needed]
DCLAD		pin clad outer diameter (mm)
DWIRE	[1.5/radial blanket 1.0]	wire diameter (mm)
HWRAP	[4.0]	S/A wrapper thickness (mm)
HCLAD	[0.4]	pin clad thickness (mm)
HWIRE	[165.0]	wire spiral repeat axial height (mm)

ITYPE	[0]	fuel pellet material type 0-MOX, 1-metal, 2-carbide, 3-nitride (not all fuel types can run all other options e.g. Minor Actinides)
TVERM	[1.98/2.0 core/radial blanket, or 0.11111, or 1.0, or 1.0]	fuel composition: metal Zr:M (w/o); O:M , C:M , N:M (a/o)
PUABND (i≤6)	[.58;.24;.14 ;.04; 0.;0.]	Pu vector: 239,240,241,242,238,Am241 (w/o)
UABND (i≤2)	[.003; .997]	U vector 235, 238 (w/o)
NOPIN	[271/radial blanket 217]	Number of pins per S/A; includes ZrH diluent pins, and tie-bar pins in ductless fuel, as well as fuel pins
DGAP	[6.0]	gap between adjacent S/As (mm)
PRING	[0.177/ radial blanket 0.109]	parameter (mm) that defines the fraction of free space (x-sectional area) within a S/A (pins considered solid) - note this is supplied in lieu of S/A a/f dimension
SMEAR	[see table]	fraction of pin internal volume (x-sectional area) occupied by pellet - note this gives any inner pellet hole diameter [ignored if DFUEL used]
THDENS	[see table]	fuel pellet material, fraction of theoretical density
VSUS (i≤5)	[.6097;.1505 ;.1950;.0251 ;.0197 - SUS316]	composition of general structural material, w/o (Fe, Cr, Ni, Mo, Mn)
ROSUS	[7.969906]	density of above material (g/cc)
VSUS2 (i≤20, j≤3)		define up to 3 'special' structural materials j=1/2/3 - pin/wrapper/wire, 5 isotopes as VSUS, plus W, then 15 dummy values (set 0?) Do not use VSUS
ROSUS2 (j≤3)		densities of up to 3 'special' structural materials (g/cc) Do not use with ROSUS & VSUS
TRRICH		fraction of transuranics (TRU) in the fuel pellet (w/o) (represents extra material added for MA burning)
TRU(i≤17)		composition (w/o) of TRU : 235U,236U, 238U,237Np,239Np,238Pu,239Pu,240Pu, 241Pu,242Pu,241Am,242Am,243Am,242Cm, 243Cm,244Cm,245Cm
ITRU		definition of TRUO2 fraction: 0 - TRUO2/(TRUO2+UO2+PUO2) 1 - TRUO2/(PUO2)
RERICH		amount of rare earth (RE) as fraction REO2/(TRUO2) (w/o)
REA(i≤22)		RE isotopic composition; numbers 36,37,...,56,57 - any not present replaced by 0, PENCIL nuclide ID's
SZRH		fractional smeared density of ZrH inside clad

VB4C		fraction of pins that are unfuelled and contain B4C cannot be used in same case as VACR
B10W		fraction (w/o) of 10B in B4C
IBASE		type of inert matrix mixed with PuO ₂ in fuel pellet: 0 - UO ₂ , 1 - MgO, 2 - CeO ₂ , 3 - BeO, 4 - Al ₂ O ₃ , 5 - ZrH ₂ , 6 - AlN
VACR		allows a fraction VACR of a 'special' material, defined in ACRND, to be mixed with the fuel. (v/o) Cannot be used with VB4C or NZRH
ACRND (i<99)		isotopic N.D. vector (PENCIL nuclide order) of 'special' material of VACR. Cannot be used with VB4C or NZRH
NZRH		No. of fuel pins that are unfuelled and contain ZrH Cannot be used with VACR
ZRH	[1.6]	composition of ZrH, ratio H:Zr
CPW	-0.97	target inner:outer peak rating ratio -ve - EOC inner : BOC outer +ve, ~1 - EOC : EOC +ve, ~100 - BOC : BOC (% value, not ratio) 0 - fixed inner:outer Pu % ratio
CKEFF	1.0048 [.996894]	target Keff for EOEC
EPW	0.002 [0.01]	allowed error on CPW
EKEFF	[0.001]	allowed error on CKEFF
PR	[0.3]	parameter to control change in Pu enrichment each iteration (%)
PTOK	[0.2]	similar to PR

[] indicates a default value.

Note - whilst values are stated as w/o (weight percent), there appears to be some confusion as to whether all such cases are actually percentages or fractions.

Additional Default Values

	MOX		Metal Fuel		Carbide Fuel		Nitride Fuel	
	core	blanket	core	blanket	core	blanket	core	blanket
SMEAR	0.876	0.908	0.750	0.800	0.850	0.900	0.850	0.900
THDENS	0.920	0.945	1.000	1.000	0.950	0.950	0.950	0.950

There are 3 different parameters that can be used for adding diluent material within fuel pins: VACR, VB4C and NZRH. Only NZRH and VB4C may be used in combination. There were errors if any 2 of these parameters were used (and whenever VB4C was used) - to avoid these make sure that the corrected load module is used

POC3A12.PENCIL52.LOAD(PENCIL).

NZRH replaces that number of fuel pins with ZrH. However, the VACR and VB4C parameters just add extra material - so

if it is to replace fuel, then the amount of fuel must be reduced (using the THDENS parameter). VACR adds material at the number densities given in ACRND (i.e. the VACR value is irrelevant). VB4C adds that fraction of the fuel volume as B₄C - it uses the fuel volume after it has been reduced using THDENS (and after any reduction for NZRH).

(Note - if VACR is used including a nuclide already in the model - fuel or steel - then you should test the program to see whether the VACR material is treated as additional material or overwrites the existing value.)

The PRING parameter is calculated as follows:

$$B = 2 \cdot (N-1) \cdot (D+d) \cdot \cos 30^\circ + D + 2 \cdot d$$

and $PRING = (E-B)/N$

where

- E hexagonal a/f distance inside wrapper
- B hexagonal a/f distance of compacted pin bundle
- N number of pin rings
- D pin diameter
- d wire diameter

Nuclide ID schemes

Nuclide	PENCIL ID	SLAROM ID	Nuclide	PENCIL ID	SLAROM ID	Nuclide	PENCIL ID	SLAROM ID
²³⁹ Pu	1	949	²³⁵ U-fp180	30	854	²⁰⁹ Bi	59	83
²⁴⁰ Pu	2	940	²⁴¹ Pu-fp180	31	814	(Mg)	60	12
²⁴¹ Pu	3	941	²⁴⁶ Cm	32	966	(Y)	61	39
²⁴² Pu	4	942	²⁴⁷ Cm	33	967	(¹⁰¹ Ru)	62	441
²⁴¹ Am	5	951	²⁴⁸ Cm	34	968	(¹⁰² Ru)	63	442
²³⁵ U	6	925	⁹⁹ Tc	35	439	(¹⁰³ Ru)	64	443
²³⁶ U	7	926	¹³⁹ La	36	579	(¹²⁹ I)	65	539
²³⁸ U	8	928	¹⁴⁰ Ce	37	580	(¹³⁵ Cs)	66	555
¹⁶ O	9	8	¹⁴² Ce	38	582	¹⁰ B	67	105
²³ Na	10	11	¹⁴⁴ Ce	39	584	¹¹ B	68	115
Fe nat.	11	26	¹⁴¹ Pr	40	591	¹⁴ N	69	147
Cr nat.	12	24	¹⁴³ Nd	41	603	H	70	1
Ni nat.	13	28	¹⁴⁴ Nd	42	604	He	71	2
Mo nat.	14	42	¹⁴⁵ Nd	43	605	Be	72	4
⁵⁵ Mn	15	25	¹⁴⁶ Nd	44	606	²⁷ Al	73	13
²³⁸ U-fp180	16	884	¹⁴⁸ Nd	45	608	(⁶ Li)	74	36
²³⁹ Pu-fp180	17	894	¹⁵⁰ Nd	46	600	(⁷ Li)	75	37
¹² C	18	6	(¹⁴⁷ Pm)	47	617	W nat.	76	74
¹⁵ N	19	157	¹⁴⁷ Sm	48	627	Ca	77	20
Zr nat.	20	40	¹⁴⁸ Sm	49	628	²³⁸ U-fp1080	78	887
²³⁷ Np	21	937	¹⁵⁰ Sm	50	620	²³⁹ Pu-fp1080	79	897
²³⁹ Np	22	939	¹⁵² Sm	51	622	²³⁵ U-fp1080	80	857
^{242m} Am	23	950	¹⁵¹ Eu	52	631	²⁴¹ Pu-fp1080	81	817
²⁴³ Am	24	953	¹⁵³ Eu	53	633	²³⁸ U-fp1800	82	889
²⁴² Cm	25	962	¹⁵⁴ Eu	54	634	²³⁹ Pu-fp1800	83	899
²⁴³ Cm	26	963	¹⁵⁴ Gd	55	644	²³⁵ U-fp1800	84	859
²⁴⁴ Cm	27	964	¹⁵⁶ Gd	56	646	²⁴¹ Pu-fp1800	85	819
²⁴⁵ Cm	28	965	¹⁵⁸ Gd	57	648			
²³⁸ Pu	29	948	Pb nat.	58	82			

Appendix II - CITATION input data

Input is in a series of blocks, the first card of which is the block identifier. There is fixed format. Where blank lines occur these must be of exactly the correct number. Blocks '002', '018', '034', '036', '091' and '093' are only used for burn-up calculations.

block '001' - control options

card 2, field 1 : 0/1 - snapshot/burn-up calculation
 fields 5-8 : +1, write macro./flux/power/source results file (to unit 31/9/32/17).
 field 19 : if >0 only uses block '008' macro. data
 card 3, options for what results are output

block '002' - burn-up control

card 2, field 1 : No. of burn-up cycles to be calculated (in PENCIL calculation, overwritten by value from geometry data)
 fields 2-4 : max. no. of burn-up timesteps per cycle (1'st, 2'nd, 3+ cycles)
 field 5 : +1/-1, do normal/normal+adjoint flux calculations at each EOC
 cards 3+4, lengths of cycles (days), reactor power MWth (power of core fraction represented)
 card 5, options for what is output during burn-up

block '003' - flux calculation definition

card 2, field 5 : geometry type (7-RZ, 13-hexZ, 14-triZ) (only used RZ, suggested 3D would be too time consuming)
 fields 8+9 : symmetry flags
 fields 11-16 : 0/1 extrapolate/reflect boundary conditions (left/top/right/bottom/front/back)
 field 18 : if >1, allows -ve flux
 card 3, convergence criteria
 card 4, field 3 : reactor power MWth (for snapshot calc.) default is 1 MW

block '004' - geometric mesh (RZ geometry)

Radial then axial mesh data. No. of meshes followed by total size of those meshes. Produces meshes of equal size, radial ones correspond to equal areas. Defines the 'volumes' where macroscopic x-sections are constant {Does not distinguish between hot & cold dimensions - the dimensions & N.D. data are used as input, without applying any expansion effects.}

block '005' - zone placement (RZ geometry)

A 'map' assigning a 'zone' number to each of the 'volumes' of data block '004'. Several 'volumes' may be in the same 'zone'. Each 'zone' number identifies a material type (i.e. different set of macro. x-section data and/or a burn-

up mesh).

block '012' - zone identification

A series of cards assigning the source of micro. x-section data for each 'zone'.

fields 1+2 : 1'st & last Nos. of consecutive set of 'zones'

field 3 : number of 'sub-zones' (i.e. batches per cycle) for this material

field 4 : position on input nuclear data file of x-section data to be used

field 6 : reference no. identifying 'zone class' ('zone class' is used in the fuel management input)

field 7 : 0/-ve for burn-up/non-depleting 'zones'

field 8 : name of 'zone' (only used in output print)

block '018' - nuclide classification

Groups nuclides into several categories, used to lump together reaction rates in output.

block '020' - nuclide Number Densities

Data is provided for all 'zones' (except in PENCIL input the core 'zones' are not included - that data is produced by the PENCIL program).

card 2, fields 1+2 : 1'st & last no's of a set of consecutive 'zones'

card 3+ pairs of nuclide identifier no. & atomic density {atoms per cc *10⁻²⁴}
[Note, no space between fields.]

block '034' - decay constants and fission yields

See CITATION manual⁽²⁾ for details. 1'st 2 cards are decay constants for 3 nuclides in core. Remaining cards are fission product yield (only necessary if standard data of micro. binary file is overwritten e.g. MA work). In each case the first card (2 values) are the 1'st and last sets of x-sections, on the micro. input file, to which burn-up is applied.

block '036' - nuclide reaction chain

See CITATION manual⁽²⁾ for details. Specifies nuclide decay & reaction relations in the core. First 2 values are (again) 1'st and last micro. datasets to which burn-up applies.

block '091' - fuel stream description

The number and arrangement of fuel streams is arbitrary (i.e. user defined). Material is removed from the reactor at the end of irradiation to a 'discharge stream', it is mixed (after any reprocessing delay) with material new to the fuel cycle from a 'feed stream' - the resultant new fuel material is placed into the reactor as a 'makeup stream'.

cards 2+ field 1 : number of fresh material 'feed streams'
field 2 : number of new fuel 'makeup streams'
field 3 : number of 'discharge streams'

field 7 : number of cycles delay between
 discharge and recycling in new fuel
 fields 9+ : identifiers of nuclides to be
 treated in fuel streams
 cards 3+ fields 1-3 : nuclide identifiers of decay parent
 & child, decay constant (sec⁻¹).

Data for each 'feed stream'

cards 4+ field 1 : 'feed stream' identification no.
 field 2 : 0/1 - use initial loading/values on
 cards 5&6, for 'feed stream' N.D.s

Data for each 'makeup stream'

card 7, field 1 : 'makeup stream' identification no.
 fields 3-5 : ID no.s of 'discharge streams' used
 field 6 : ID no. of 'feed stream' used
 field 9 : fraction of material available from
 'discharge streams' to be used
 cards 8+ for each nuclide on card 2, fraction of
 'recovered' material in 'discharge streams' to be
 used

Data for each 'discharge stream'

card 9, field 1 : 'discharge stream' identification no.
 cards 10+ for each nuclide on card 2, fraction of
 discharged material 'recovered' during
 reprocessing

block '093' - fuel management

card 2, field 2 : {uncertain of meaning, CITATION manual
 description is unclear}
 field 4 : no. of fuel cycles calculated
 cards 6+ Each card specifies one 'fuel movement'
 field 1 : 'identification no. of 'fuel movement'
 field 2 : 'feed stream' used for new fuel before
 recycled material is available
 (0 - 'closed loop', no refuelling)
 field 3 : 'feed stream' used in mass balance
 calculation
 field 4 : 'makeup stream' used for new fuel
 after recycled material is available
 field 8 : 'discharge stream', destination of
 removed fuel
 field 10-12 : cycle no. when this 'fuel
 movement' ends (if <>0)/starts/
 frequency of use
 fields 13+ : 'zone class' to which 'fuel
 movement' is applied, if >1 value
 is a shuffling through 'zones'
 with defuelling from the last one
 (a closed loop, no defuelling, is
 possible); with 'sub-zones' only
 1 'sub-zone' is moved each cycle,
 in order {CITATION manual is
 confused over 'zones' and 'zone
 classes'}

block '999' - end of data marker

Appendix III - PERKY input data

Line 1 'PERKY' (program ID)
 Line 2 1'st fuel material region; 'SLAROM'
 (creating program ID)

Title Line
 Block '001' - general data
 Line 2
 field 2 no. of energy groups
 field 4 no. of material regions input (max 40)
 field 5 1/0 - RZ/XY or XYZ geometry
 field 6 1 if an exact perturbation calc.
 (else 0)

Block '002' - microscopic data input
 Line 3
 field 1 total no. of microscopic datasets
 input
 field 2 no. of microscopic datasets input by
 JOINT
 field 3 no. of microscopic datasets input by
 2'nd input file (I have not used this
 option)
 field 4 no. of microscopic datasets input
 within this file (I have not used this
 option)
 field 5+ id. no. assigned to each microscopic
 dataset

For each material region for which JOINT microscopic
 data is input:
 line 1 microscopic dataset ID no.; no. of
 isotopes with microscopic data input
 (max. 20)
 line 2 name of material region providing
 microscopic data; 'SLAROM' (creating
 program ID)
 line 3 ID no.s of nuclides for which
 microscopic data is input

Last line Pairs of: material region ID no.;
 microscopic data ID no. of corresponding
 data (blank where no such data is input)

Block '003' - macroscopic data input
 Line 3+
 field 1 total no. of macroscopic datasets
 input
 field 2 no. of macroscopic datasets input by
 JOINT
 field 3 no. of macroscopic datasets input by
 2'nd input file (I have not used this
 option)
 field 4 no. of macroscopic datasets input
 within this file (I have not used this
 option)
 field 5+ material region ID no. assigned to
 each macroscopic dataset

Line 4 material region supplying fission spectrum
 data; 'SLAROM' (creating program ID)

Line 5
 field 1 no. of energy groups
 field 2 no. of downscatter groups (max 30?)
 field 3 no. of upscatter groups

Line 6+ for each material region providing JOINT
 macroscopic data input: material region
 name; 'SLAROM' (creating program ID)

Block '004' - calculation type
 Line 2 1 - delayed neutron fraction
 2 - prompt neutron lifetime
 3 - reactivity worth map
 4 - 1'st order perturbation
 5 - exact perturbation

Block '005' - delayed neutron data
 Line 2
 field 1 no. of energy groups
 field 2 no. of delayed neutron precursor
 families
 field 3 no. of nuclides giving rise to delayed
 neutrons
 field 4+ ID no. of each delayed neutron
 producing nuclide

Line 3 no. of energy groups
 For each delayed neutron producing nuclide:
 Line 1 nuclide ID; no. of delayed neutron
 precursor families
 For each delayed neutron precursor family
 fraction of neutrons that are delayed (by
 energy group)
 energy spectrum of delayed neutrons

Block '006' - neutron lifetime data
 Line 2+ upper bound lethargy of each energy group

Block '007' - fission distribution data (I have not used
 it)

Block '008' - perturbation data
 Line 2
 field 1 0/1 - microscopic/macrosopic data
 used to define perturbed state
 field 2 1 for 'by nuclide' perturbation,
 0 otherwise {meaning not certain}
 field 4 no. of material regions where
 perturbation changes occur
 field 5+ ID no.s of material regions to be
 changed

Line 3 name of 1'st perturbed material region;
 'SLAROM'

'by nuclide' case
 Line 4
 field 1 no. of perturbed nuclides (max 10)
 field 2+ ID no.s of perturbed nuclides

Line 5 ID no.s of burnable isotopes included in
 flux calculation (max 10) {uncertain of
 meaning}

For each material region with perturbation changes
 line 1
 field 1 ID no. of normal microscopic data
 field 2 ID no. of perturbed microscopic
 data

field 3 no. of perturbed nuclides (max 10)
For each perturbed isotope
field 1 ID no. of perturbed nuclide
other cases
Line 4 pairs of microscopic data or material
region
ID no.s: a normal dataset, followed by the
perturbed version that replaces it
Block '999' - end of data marker.
Followed by one blank line.

Note that the ID Nos. given to unperturbed microscopic datasets and material regions should correspond to the zone numbers in the CITATION calculation that produced the unperturbed flux file used in the PERKY calculation.

Note that the input of microscopic and macroscopic datasets via JOINT, and of delayed neutron data, is not as described in the (Japanese) PERKY manual⁽⁵⁾.

Appendix IV - MOSES input data

The manual for MOSES is reference 9. It gives a fuller description of the input; unfortunately, it is only available in Japanese.

There are 2 input file names to be set to run a case -
 FT03F001 is a microscopic nuclear data file, in CITATION (binary file) format
 FT05F001 is the case input data.
 Other input files, which I have not used, are -
 FT08F001 is used for debugging options
 FT69F001 is fuel shuffling data in refuelling sequences

If a case is large, then it may be necessary to alter the size of temporary workfiles used by the program (these have to be user specified!). There are a series of parameters such as N1, N2, NX etc. specified in the JCL - if a job fails due to insufficient space, look for the part of the output in the failed job where recommended values are given for these parameters.

Input is in a series of blocks, in order, each identified by a **KEYWORD** at its beginning. Some blocks are ended by the words **EOI** or **EOB**. Input is in free format.

- Title** Starts with 2 title lines (no 'TITLE' keyword).
- CONTROL** Calculation type - Steady State or Burn-up.
 6 values (0/1). Set just the first 3 values to 1 for a diffusion calculation, all 6 for burn-up.
- EDIT** Output option control.
 4 sets (lines) of integer values.
 [Details not known.]
- EDITD** Detailed S/A output control.
 3 sets each of 50 values. No.s of S/As for which detailed output is produced; for diffusion, burn-up and N.D. calculations.
 Will give data for peak S/As anyway (with standard **EDIT** values).
- EOB**
- LOADING** Core loading/refuelling step selector.
 The ID number of a core S/A map dataset from **REFMAP**. The S/As specified by this dataset will be loaded, unloaded or both, as specified by **REFUEL** (can be e.g. full core description or a refuelling step). (0 for a calculation step with no change in core layout.)
- GEOM** Basic geometric data -
NSYM symmetry - no. of 60° sectors modelled (1,2,3 or 6); the symmetry is rotational.

NLAY no. of S/A rings in model
 NPLN no. of axial planes (mesh centred flux)
 (supposedly limited to 40, but have used larger
 number without problems)
 NAZN no. of axial material regions. Defines where
 axial boundaries in material compositions are
 allowed. {supposedly max 14, but I accidentally
 used 15, without any apparent error}.
 NGRP no. of energy groups
 PITCH hexagonal mesh a/f pitch (cm)
 MESHZ no. of uniform flux meshes in each axial material
 region (NAZN values, adding up to NPLN).
 DELTZ height (cm) of each axial region (NAZN values)

CPARA1

Calculation type and time step control -
 NMES 1 - 1 hex mesh per S/A;
 2 - 1 hex mesh per S/A, interpolated to 6 rating
 (not n.d.) meshes per S/A (Askew
 method);
 3 - 6 meshes per S/A.
 NCAL 0 - direct flux calculation;
 1 - direct & adjoint fluxes calculations;
 2 - source calculation.
 NITR/NIIT no. of outer/inner iterations. (For NMES=3, try
 increasing NIIT from the default 3 to 8.)
 NTIM flux solution curtailed after this many seconds
 CPU time
 8 values (not to be used)
 DELDAY burn-up time (days) for this calculation step
 NSMALL number of sub-steps a burn-up calculation step is
 divided into (flux is re-evaluated each sub-step)

CPARA2

Calculation convergence control -
 EPSRMD Keff convergence limit (have reduced to 5.0e-6)
 EPSF flux convergence limit (have reduced to 1.0e-4)
 EPSMIN/EPSSMAX extreme Keff value bounds (use default
 values, 0.5/1.5)
 METHOD burn-up calculation method.
 (Option 1 always used.)
 1 - 'matrix exponential'
 2 - 'average generation rate'.
 8 values (not to be used, only for METHOD=2)

EOB

AXIAL

S/A axial composition data.
 Defines an axial map of material types onto axial
 regions; boundaries only allowed where already
 set by DELTZ. These regions also define the
 axial n.d. mesh in a burn-up calculation. For
 each S/A type to be used in the model (max 40) -
 line 1 NREGZ, followed by NREGZ values of material type
 ID
 line 2 NREGZ values of axial region height (sum equals
 sum of DELTZ values)
 End with **EOI** marker.

{In earlier calculations, I had problems when I used

NREGZ=14 axial zones in a model with NAZN=14. This was a program bug, now corrected - errors were apparent when checking the axial map of S/A materials in the MOSES output.}

ASYGROUP S/A output data.

For each of the S/A types of **AXIAL**, and in the same order, the following items -

- field 1 1 - fixed; 2 - axially moving (an absorber rod). 'Axially moving' allows the use of **CRPOS** to reposition rod insertions between calculation steps
- field 2 an ID number (max 15) used in the MOSMAP core map
- field 3 0 or 1 (always use 1)
- field 4 1 - inner core; 2 - outer core; 3 - radial blanket; 4 - shield, rods etc.
Used in print output.
- field 5 if -ve then label 'field 6', rather than this value, is used in plotted output
- field 6 S/A type ID label
- field 7 S/A type name (used in output) - 2 separate words
End with **EOI** marker.

REGION Material type data.

For each of the material types (max 60) used in the S/A **AXIAL** mapping, the following -

- field 1 position of entry on nuclear data file to be used for microscopic nuclear data
- field 2 material type ID - used to define materials in some of the output, allows several to be combined
- field 3 flag used to set printing option (*always use 1*)
- field 4 1 - fuel; 2 - axial breeder; 3 - axial shield; 4 - radial breeder; 5 - radial shield; 6 - rods
- field 5 flag used to set plot option (*always use 1*)
- field 6 material type identifying label
- field 7 material type name (used in output) -
2 separate words
End with **EOI** marker.

MICROXS Nuclear data (general).

- NLIB source of nuclear data (always use 1 - FT03)
- NSET no. of nuclear data entries in source file
- NNUC no. of nuclides used in the MOSES calculation
- NRAC no. of entries in nuclear dataset that include burnable nuclides (must these be the 1'st entries?)
- NHAT thermal energy option: 0 - only fission x-sections 1 - capture & fission (2 - unknown, do not use)
- NFGP no. of groups > 0.1 Mev (used in printout only)
- CHI spectrum of neutrons produced by fission

XSNUC Nuclear data, for each of the NNUC nuclides -

- IDNUC the nuclide ID number used in the nuclear dataset (i.e. PENCIL ID no.)
- NAMNUC nuclide name, used in print output
- IFIS fissile nature identifier: -1 - fission products; +1 - uranium; +2 - plutonium (inc. Am241);

0 - others. Those set +1/+2 will have fission reactions included in the burn-up calculations. Defines materials for breeding ratio calculation. nuclide classification: 1 - fissile; 2 - fertile; 3 - other fuel nuclides; 4 - fission products; 5 - oxygen; 6 - coolant & structure; 7 - B4C; 8 - 'anything else'. Used to define materials for fuel cycle balance. MA are type 3, diluent type 6.

NBURN always use 0

NANAL if 1, use standard values of ATW & EFISS (rather than data input below); if 2 also use standard

KLIB DECAF values

ATW atomic weight

EFISS energy per fission

ECAPT energy per capture

DECAF decay time constant

REACT1 Decay chain data. A series of lines with 4 values followed by a card with '0'. Max. no. of entries not specified; if too many are used, unflagged array bound errors can cause subtle errors (e.g. in Keff); appears OK with 19 entries. Was initially advised to include fission products producing themselves by neutron capture - seems dubious, since other isotopes are not needed here to make capture effective (not in MOSES manual example).

field 1 ID of nuclide undergoing reaction (IDNUC value)

field 2 ID of nuclide produced by reaction

field 3 reaction type: 1 - beta decay; 2 - (n,gamma); 3 - (n,alpha); 4 - (n,p); 5 - (n,2n); 6 - (n,d); 7 - (n,t); 8 - special, used when x-section library excludes the (n,2n) reaction (takes value 'field 4' x fission x-section).

field 4 fraction of reactions that are of this type, where more than one reaction product is possible

YIELD Fission product yield data. A series of lines with 3 values on, all nuclides with IFIS > 0 must be included. As with **REACT1**, too many values cause array bound errors; it appears to work with up to 14 values, but am advised to use no more than 10 values.

IFP ID of fission product nuclide (IDNUC value)

IFS ID of a nuclide yielding fission product IFP

YIELD fraction of IFS that end up as IFP

ATDEN Number density data. For each of the material types in **REGION**, and in the same order - A list of number densities (atom/cc * 10⁻²⁴) for each of the nuclides in **XSNUC**, in the same order. End with **EOI** marker.

BOUNDARY Boundary condition data. One value for each energy group, for the radial, then upper, then lower boundaries. Default value

is 0.4692; normally use vacuum boundary, 0.0; reflective boundary is 1.0.

REFUEL Burn-up calculation fuel cycle balance data. Each line is associated with one set of input under **REFMAP**. Exactly 10 lines of input -
 field 1 1/0 indicates fuel is/is not loaded
 field 2 1/0 indicates fuel is/is not removed
 field 3 refuelling step length (days) (not necessarily the same as calculation step DELDAY).

REFMAP Core loading/fuel movement S/A sets. Identifies S/As that are loaded and/or unloaded, according to **REFUEL** data, when a **REFMAP** set is invoked by **LOADING**. One set of data for each non-zero line under **REFUEL** (max 10 sets). Each set has a title line, followed by lines identifying a series of S/As -
 I1 first of a series of continually numbered S/As
 I2 last of a series of continually numbered S/As
 K ID of S/A type assigned to positions I1 to I2 (For ease of input, later entries can overwrite S/A type IDs assigned in an earlier line.)
 - there is final line containing values 0 0 0.
 End with **EOI** marker.

PLANT
 CPOWER reactor power (MWth). For NSYM<>6 this power must be reduced proportionately.
 TIN reactor inlet temperature
 [not used in calculation]

CRPOS A series of sets of 40 values (1 for each S/A type allowed). Provided ASYGROUP(1)=2, indicates how far the rod positions are moved from those set in **AXIAL**. +ve values indicate upward movement; all material boundaries move, the top/bottom material is used to fill in any space created. Is not restricted to mesh positions defined in **GEOM**.

DBURNUP [Optional]
 Data for option to interpolate 1 mesh per S/A n.d. representation to 6 meshes per S/A, in selected S/As. OPTION NOT YET IMPLEMENTED IN MOSES.

ASMRC and FIXSRC [Optional]
 Neutron source data (I have not used this option).

EOB

Multiple cases, or a series of steps within a single burn-up calculation, are input by repeating the input data, including the TITLE and all EOB lines. It is generally only necessary to input keywords where values change,

though all values for that keyword have to be input. It appears likely that at least one keyword should be included for each **EOB** block - **CONTROL**, **LOADING** and **CRPOS** are the favourite candidates for permanent inclusion in the 3 blocks.

It should be possible to use either **CRPOS** or **REFMAP** to change rod insertions. **REFMAP** has the advantage that it does not move any detailed above/below core structure, all that is required is to have S/As for the different rod insertions defined under **AXIAL**, and to "refuel" the rod S/As.

The mesh representation gives values for flux, rating etc. at the centres of meshes. This even applies to the extrapolated 6 point values obtained using the Askew type calculation. (A proposed future version will produce values that are actual 'peak interpolated' values, rather than centre of triangle values.)

Sub-assembly numbering starts with 1 at the core centre, then each succeeding ring is numbered in order, clockwise. The initial S/As of each ring, together form a straight line of S/As, each sharing edges with its neighbours. There is an interactive program **MOSMAP**, which will take any **MOSES** input dataset and produce plots of the S/A layout (useful for data checking).

The **MOSMAP** command file - **POC3C##.CMDPROC.CLIST(MOSMAP)** - is listed in my Macro enabling file - **.TSS.CLIST** - which is invoked as described under the **RZOUT3** calculation: at the command prompt input

```
STREAM COMMAND
EX TSS
LOGON
```

To start **MOSMAP**, at the command prompt just type **MOSMAP** - you are prompted for the data file name, this must be the full name path (since the **MOSMAP** procedure file is owned by another user). The output goes to a printer situated in an adjacent building (the building where, at present, Aoyagi-san works 3 days per week).

Appendix V - Core Expansion Reactivity Effect

The following analysis is that used for the Pu burner core⁽⁹⁾, it was in turn based on that used for the PNC design study on a 600 MW(e) class reactor⁽¹⁰⁾.

Two types of computer calculation are done to produce components of the thermal expansion effects on reactivity.

One calculation type is 1'st order PERKY calculations. These calculate the reactivity effects of density changes in the fuel, coolant and structural materials:

$$K_F = \frac{\Delta k / kk'}{\Delta \rho_F / \rho_F} \quad , \quad K_C = \frac{\Delta k / kk'}{\Delta \rho_C / \rho_C} \quad , \quad K_S = \frac{\Delta k / kk'}{\Delta \rho_S / \rho_S}$$

(Note that any diluent material goes in the structural category.)

The second type of calculation is CITATION calculations for the core, but with either axial or radial mesh sizes slightly increased (say 1%), the number densities remain the same. These calculate the reactivity effect of changes in core height and radius:

$$K_H = \frac{\Delta k / kk'}{\Delta H / H} \quad , \quad K_R = \frac{\Delta k / kk'}{\Delta R / R}$$

The thermal expansion effects have to be added to temperature effects in cross-sections to get the complete reactivity effect of changing reactor temperatures. The following shows how the core expansion reactivity effect is calculated from the above factors, which are produced by the 'shutdown margin' calculation block.

The various core components undergo different amounts of thermal expansion, the overall effect is obtained from summing the following 5 factors:

$$F_1 = \Delta T_{sup} \cdot \alpha_{steel}(T_{sup}) \cdot \left(K_R - 2 \cdot (K_F + K_S) + 2 \cdot \frac{(1 - V_{Na})}{V_{Na}} \cdot K_C \right)$$

$$F_2 = - \Delta T_{wrap} \cdot \alpha_{steel}(T_{wrap}) \cdot 2 \cdot \frac{V_{wrap}}{V_{Na}} \cdot K_C - \Delta T_{clad} \cdot \alpha_{steel}(T_{clad}) \cdot 2 \cdot \frac{V_{pin}}{V_{Na}} \cdot K_C$$

$$F_3 = \Delta T_{fuel} \cdot \alpha_{fuel}(T_{fuel}) \cdot (K_H - K_F)$$

$$F_4 = - \Delta T_{clad} \cdot \alpha_{steel}(T_{clad}) \cdot \frac{V_{clad}}{(V_{clad} + V_{wrap})} \cdot K_S - \Delta T_{wrap} \cdot \alpha_{steel}(T_{wrap}) \cdot \frac{V_{wrap}}{(V_{clad} + V_{wrap})} \cdot K_S$$

$$F_5 = \Delta T_{Na} \cdot \alpha_{Na}(T_{Na}) \cdot K_C$$

V - volume fraction

α - material thermal expansion coefficient, temp. dependent (density coefficient for Na coolant)

ΔT - temperature change
sup - support diagrid
wrap - wrapper
pin - all pins (including contents)

The support diagrid effects overall radial expansion (F1) - core radius increases, this also causes solid material densities to reduce; coolant fills the added volume. (Radial expansion is in 2 dimensions, hence the factor 2.)

Radial expansion within the S/A displaces some coolant (F2) - wrapper volume and pin volume (fuel & diluent) expand (pin expansion assumed driven by steel clad, not pin contents).

Fuel expansion effects overall axial expansion (F3) - assumed not stuck to clad, core volume increases, but fuel density decreases.

Axial expansion of non-fuel components (F4) - clad, wrapper and diluent densities reduce; diluent is assumed to expand with the clad. (Diluent set partly to each of the clad & wrapper temperatures - necessary to make equation tractable.)

Coolant density changes (F5).

As well as the calculations of PERKY and CITATION to produce reactivity/temperature coefficients, need to have values of expansion coefficients and of temperatures appropriate to the various materials. The following values were used for the Pu burner core⁽⁹⁾, they were taken from the PNC design study on a 600 MW(e) class fast reactor, it may be appropriate to calculate replacement values -

Operating temperatures (°C)	
support	380
coolant	455
wrapper	455
clad	475
fuel	1073
Cold shutdown temperature	200°C
Linear expansion coefficients ($\Delta l/l$ per °C)	
fuel	1.439×10^{-5}
clad	2.016×10^{-5}
wrapper	2.003×10^{-5}
support	1.949×10^{-5}
Density expansion coefficient ($\Delta \rho/\rho$ per °C)	
coolant	-2.841×10^{-4}

Appendix VI - Shutdown Margin Components

The calculations of the 'rod worth' block provide nominal values of the rod group worth for both the main (PCR) and backup (BCR) rod groups, from a difference in Keff values. To translate these to values of shutdown margins, it is necessary to apply allowances for uncertainties in the calculated rod worth values, and to compare the result with the shutdown requirements. The following lists the various components of rod worth uncertainty and shutdown requirements that are included in the PNC calculation route; the values used for the Pu burner core⁽⁹⁾ are given below, they were taken from the PNC design study on a 600 MW(e) class fast reactor⁽¹⁰⁾ - it may be appropriate for you to obtain a different set of values.

Components of Rod Worth Uncertainty

Effect Adjusted	Factor
Effect of Condensation to 7 energy groups	x0.972
Combined Mesh and Transport correction factor	x1.021
Control Rod Heterogeneity correction factor	x0.88
Rod ¹⁰ B Burn-up correction factor	x0.95
Difference between MOSES program and reference (CITATION) used in assessing corrections	x1.0
Adjustment for difference from experimental study (based on JUPITER core ZPPR-10C)	x1.018
Allowance for calculational uncertainty (2 σ)	x0.9
Total Correction Factor	x0.761

Components of Shutdown Requirement

Components of Main Rod Group (PCR) Shutdown Requirement	
Effect of reducing reactor temperature to 200°C	α
Reactivity loss per cycle	β
Margin for control at EOC	0.2 % $\Delta k/kk'$
Margin for variations in refuelling batches	0.1 % $\Delta k/kk'$
Uncertainties in cycle length #	
Reactor measured power uncertainty (2 σ)	0.24 % $\Delta k/kk'$
Fuel fabrication uncertainties (2 σ)	0.3 % $\Delta k/kk'$
Uncertainty in reactivity loss/cycle (2 σ)	20% of β
Uncertainties in shutdown margin #	
Uncertainty in calculated Keff (1 σ)	0.44 % $\Delta k/kk'$
Fuel fabrication uncertainties (2 σ)	0.3 % $\Delta k/kk'$
Components of Backup Rod Group (BCR) Shutdown Requirement	
Effect of reducing reactor temperature to 200°C	α
Margin for reactivity inserted during fault	0.2 % $\Delta k/kk'$
Uncertainty, based on JOYO measurements (2 σ)	0.24 % $\Delta k/kk'$

components are combined in quadrature

The values of α and β are assessed for each use of the calculational route.

α is the combined reactivity worths of the core expansion (see previous addendum) and changing nuclear data.

β is obtained from the S-J-C calculations of the 'basic fuel cycle' block.

Appendix VII - Example Datasets

The following pages list the various JCL and data files referred to in Section 6 of this note. All the files should be available on the FACOM computer, under uid POC0B17, if not immediately accessible then they are all available from my archive tape cartridges.

In the listings there are various line number fields given at the end (sometimes the beginning) of the lines - these can be ignored, they are an artifact of the FACOM file storage and editing system.

The following is a list of all the filenames, with the page number and the ID no. used in Section 6.

JCL file [1]	.OPJCL.CNTL(GDI1PEN)	83
JCL file [2]	.OPJCL.CNTL(GDI1CDEN)	84
JCL file [3]	.TSS.CLIST	85
JCL file [4]	.TSSMAC.CLIST(RZOUT3)	85
JCL file [5]	.OPJCL.CNTL(GDG0SJC)	85
JCL file [6]	.OPJCL.CNTL(GDI1MASS)	87
JCL file [7]	.OPJCL.CNTL(GDI1SC)	88
JCL file [8]	.OPJCL.CNTL(GDI3PSLA)	89
JCL file [9]	.OPJCL.CNTL(GDI3P70N)	90
JCL file [10]	.OPJCL.CNTL(GDI3P70V)	91
JCL file [11]	.OPJCL.CNTL(GDI3P18N)	92
JCL file [12]	.OPJCL.CNTL(GDI3P70D)	93
JCL file [13]	.OPJCL.CNTL(GDI3PNL)	94
JCL file [14]	.OPJCL.CNTL(GDI3PDF)	95
JCL file [15]	.OPJCL.CNTL(GDI3PEV)	96
JCL file [16]	.OPJCL.CNTL(GDI3PFD0)	97
JCL file [17]	.OPJCL.CNTL(GDI3PFD1)	98
JCL file [18]	.OPEDID.DATA(DUMPGDI1)	98
JCL file [19]	.OPJCL.CNTL(GDI1MY)	102
JCL file [20]	.OPJCL.CNTL(GDI1XSLA)	103
JCL file [21]	.OPJCL.CNTL(GDI1X70N)	104
JCL file [22]	.OPJCL.CNTL(GDI1X70D)	106
JCL file [23]	.OPJCL.CNTL(GDI1PERK)	107
JCL file [24]	.OPJCL.CNTL(GDI1MBRN)	110
Data file [1]	.OPDEND.DATA(GDI1P)	112
Data file [2]	.OPCITD.DATA(O975M165)	112
Data file [3]	.OPEDID.DATA(GDI1RZ)	116
Data file [4]	.OPFORT.DATA(GDI1RZ)	116
Data file [5]	.OPDEND.DATA(GDG0S)	118
Data file [6]	.OPCITD.DATA(PHA1N70)	120
Data file [7]	.OPEDID.DATA(PFC1F)	120

Data file [8]	.OPEDID.DATA(PFC1C)	121
Data file [9]	.OPEDID.DATA(PFC1M)	121
Data file [10]	.OPDEND.DATA(GDI1S)	122
Data file [11]	.OPCITD.DATA(PEE1D700)	124
Data file [12]	.OPDEND.DATA(GDI3PRT)	126
Data file [13]	.OPCITD.DATA(PIJ3P70N)	129
Data file [14]	.OPEDID.DATA(PIJ3PF)	130
Data file [15]	.OPEDID.DATA(PIJ3PC)	130
Data file [16]	.OPCITD.DATA(PIJ3PV70)	131
Data file [17]	.OPCITD.DATA(PIJ3PADJ)	132
Data file [18]	.OPCITD.DATA(PIJ3P70D)	133
Data file [19]	.OPEDID.DATA(PIJ3PNL)	133
Data file [20]	.OPEDID.DATA(PIJ3PDNX)	134
Data file [21]	.OPEDID.DATA(PIJ3PEV)	135
Data file [22]	.OPEDID.DATA(PIJ3PFD0)	136
Data file [23]	.OPEDID.DATA(GDI3PFD1)	136
Data file [24]	.OPMOSD.DATA(GDI1DY)	140
Data file [25]	.OPDEND.DATA(GDI1RZEC)	143
Data file [26]	.OPCITD.DATA(PFC3N70X)	145
Data file [27]	.OPEDID.DATA(PFC3FX)	145
Data file [28]	.OPEDID.DATA(PFC3C18X)	146
Data file [29]	.OPEDID.DATA(PFC3MX)	146
Data file [30]	.OPEDID.DATA(NAMEF#XX)	147
Data file [31]	.OPEDID.DATA(IDNUMXX)	147
Data file [32]	.OPEDID.DATA(CHIMG)	149
Data file [33]	.OPEDID.DATA(ZONE#XX)	149
Data file [34]	.OPMOSD.DATA(GDI1MBRN)	149

If the above files are no longer present in the permanent filespace under uid POC0B17, they can be retrieved from archive on my magnetic cartridge, identifier SNHBU3.

The JCL file to run the retrieval is .ARCHTAPE.CNTL(REST30), a listing is included after the last data file. It may be necessary to alter the names of the files that are retrieved (e.g. to temporary filespace), do this by altering the SYSUT2 data in each step of the JCL.

Example JCL Files

JCL file [1] .OPJCL.CNTL(GDI1PEN)

PENCIL, 'iterative condensation' & 'basic fuel cycle'

```

//POCOB17D JOB (),PNCL850R,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010022
// CLASS=C,TIME=0030 00020020
//----- 00030007
//IMAGELIB DD DSN=POC3A12.IMAGELIB,DISP=SHR 00040007
//DATAP1 EXEC PGM=DATAP1 00090007
//PSYS DD SUBSYS=(VPCS) 00100017
//STEPLIB DD DSN=POC3AA3.LIB.LOAD,DISP=SHR,LABEL=(, , , IN) 00110000
//* ++++++ INPUT GEOMETRY DATA 00120010
//FT05F001 DD DSN=POCOB17.OPDEND.DATA(GDI1P),DISP=SHR,LABEL=(, , , IN) 00121021
//* ----- 00122010
//FT06F001 DD SYSOUT=* 00130000
//----- 00140000
//* PENCIL 00150000
//----- 00160000
//PENCIL EXEC PGM=PENCIL,COND=(4,LT) 00170007
//PSYS DD SUBSYS=(VPCS) 00180017
//STEPLIB DD DSN=POC3A12.PENCILS2.LOAD,DISP=SHR,LABEL=(, , , IN) 00190043
//* CORRECTED PENCIL VERSION - VB4C & NZRH COMPATIBLE 00200043
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(20,2)), 00210000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00220000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00230000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00240000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00250000
//FT05F001 DD DSN=&&CIT,DISP=(NEW,PASS),UNIT=WORK, 00260007
// SPACE=(TRK,(10,2)),DCB=(RECFM=FB,LRECL=80,BLKSIZE=26720) 00270007
//FT06F001 DD SYSOUT=*,DCB=(RECFM=UA,BLKSIZE=137) 00290000
//FT07F001 DD UNIT=SYSDA,SPACE=(TRK,(50,50)),DCB=*.FT01F001 00300000
//* ++++++ 7 GROUP MICRO DATA 00310010
//FT08F001 DD DSN=POCOB17.OBH2HBC3.MIC18G3,DISP=SHR,LABEL=(, , , IN) 00311017
//* ----- 00312010
//*---( OUTPUT FLUX MAP )-----** 00320000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00330000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00340000
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00350000
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00360000
//*---( RESTART FILE )-----** 00370000
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00380000
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00390000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00400000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00410000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00420000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00430000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00440000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00450000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00460000

//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00470000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00480000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00490000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00500000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00510000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00520000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00530000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00540000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00550000
//*---( MACRO CROSS SECTIONS )----- 00560000
//FT31F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00570000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00580000
//FT33F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00590000
//*---( OUTPUT RESULT FILE )----- 00600000
//FT34F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00610000
//FT35F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00620000
//FT36F001 DD DUMMY 00630000
//FT37F001 DD DUMMY 00640000
//*---( OUTPUT LIST )----- 00650000
//FT51F001 DD DUMMY 00660000
//FT52F001 DD DUMMY 00670000
//FT80F001 DD DUMMY 00680000
//FT81F001 DD DUMMY 00690000
//*---( OUTPUT FLUX FILE )----- 00700000
//FT91F001 DD DSN=&&FLUX,DISP=(NEW,PASS),SPACE=(TRK,(20,10),RLSE), 00710007
// UNIT=WORK,DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00720007
//*---( OUTPUT POWER FILE )----- 00730000
//FT92F001 DD DSN=&&POWER,DISP=(NEW,PASS),SPACE=(TRK,(1,1),RLSE), 00740007
// UNIT=WORK,DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00750007
//* ++++++ INPUT GEOMETRY DATA 00770011
//FT93F001 DD DSN=POCOB17.OPDEND.DATA(GDI1P),DISP=SHR,LABEL=(, , , IN) 00771021
//* ----- 00772010
//* ++++++ INPUT CITATION FILE 00790010
//FT94F001 DD DSN=POCOB17.OPCITD.DATA(O975M165),DISP=SHR,LABEL=(, , , IN) 00791010
//* ----- 00792010
//* ++++++ OUTPUT TABLE DATA LIST 00810010
//FT95F001 DD DSN=POCOB17.PQINT.DATA(TABLE),DISP=SHR 00811015
//* ----- 00812010
//* ++++++ OUTPUT APPLE DATA FILE 00830010
//FT96F001 DD DSN=POCOB17.PQINT.DATA(APPLE),DISP=SHR 00831015
//* ----- 00832010
//* ----- 00840000
//* PENEDIT 00850000
//*----- 00860000
//CITCPY EXEC PGM=PENEDIT 00870007
//PSYS DD SUBSYS=(VPCS) 00880017
//STEPLIB DD DSN=POC3A12.PENEDIT.LOAD,DISP=SHR,LABEL=(, , , IN) 00890000
//FT02F001 DD DSN=&&CIT,DISP=(OLD,DELETE),UNIT=SYSDA 00900000
//* ++++++ OUTPUT CITATION DATA 00910010
//FT12F001 DD DSN=POCOB17.PQINT.DATA(CIT),DISP=SHR 00911015
//* ----- 00912010
//FT06F001 DD SYSOUT=* 00920000
//*----- 00930000

```

```

/** APPLE4 00940000
----- 00950000
//APPLE4 EXEC PGM=APPLE4 00960007
/**PSYS DD SUBSYS=(VPCS) 00970017
//STEPLIB DD DSN=POC3AA3.APPLEW.LOAD,DISP=SHR,LABEL=(, , , IN) 00980000
//FT01F001 DD DSN=&&AW1,DISP=(NEW,PASS),UNIT=SYSDA, 00990000
SPACE=(TRK,(10,2)),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01000000
//FT02F001 DD DSN=&&AW2,DISP=(NEW,PASS),UNIT=SYSDA, 01010000
SPACE=(TRK,(10,2)),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01020000
//FT03F001 DD SYSOUT=*,DCB=(RECFM=UA,BLKSIZE=133) 01030000
/**
***** OUTPUT APPLE FILE01040010
//FT05F001 DD DSN=POCOB17.PQINT.DATA(APPLE),DISP=SHR,LABEL=(, , , IN) 01041015
----- 01042010
//FT06F001 DD SYSOUT=* 01050000
//FT07F001 DD DUMMY 01060000
//PLOTLOG DD SYSOUT=* 01070000
//PLOTPRM DD * 01080000
SCALE=0.697 01090000
/* 01100000
//GDFILE DD SYSOUT=* 01110000
----- 01120000
/** TABLE1 01130000
----- 01140000
//TABLE EXEC PGM=TABLE1 01150000
/**PSYS DD SUBSYS=(VPCS) 01160017
//STEPLIB DD DSN=POC3A12.LIB.LOAD,DISP=SHR 01170000
***** OUTPUT CITATION DATA01180010
//FT01F001 DD DSN=POCOB17.PQINT.DATA(CIT),DISP=SHR 01181015
----- 01182010
//FT02F001 DD DUMMY 01190000
//FT03F001 DD DUMMY 01200000
/**
***** OUTPUT APPLE FILE01210010
//FT04F001 DD DSN=POCOB17.PQINT.DATA(APPLE),DISP=SHR,LABEL=(, , , IN) 01211015
----- 01212010
/**
***** OUTPUT TABLE LIST01213010
//FT05F001 DD DSN=POCOB17.PQINT.DATA(TABLE),DISP=SHR,LABEL=(, , , IN) 01220015
----- 01221010
/**
***** INPUT 7 GROUP MICRO DATA01222010
//FT07F001 DD DSN=POCOB17.OBHZHBC3.MIC18G3,DISP=SHR,LABEL=(, , , IN) 01230017
----- 01231010
/**
***** INPUT GEOMETRY DATA01232011
//FT08F001 DD DSN=POCOB17.OPDEND.DATA(GDI1P),DISP=SHR,LABEL=(, , , IN) 01240021
----- 01241010
//FT06F001 DD SYSOUT=*,FLASH=TBL1 01250000
----- 01260000
/** RZPLOT 01270000
----- 01280000
//RZPLOT EXEC PGM=RZPLOT,COND=(4,LT) 01290007
/**PSYS DD SUBSYS=(VPCS) 01300017
//STEPLIB DD DSN=POC3AA3.LIB.LOAD,DISP=SHR 01310000
***** OUTPUT CITATION DATA01320010
//FT05F001 DD DSN=POCOB17.PQINT.DATA(CIT),DISP=SHR,LABEL=(, , , IN) 01321015
----- 01322010
//FT06F001 DD SYSOUT=* 01330000
//GDFILE DD SYSOUT=* 01340000
//PLOTPRM DD DUMMY 01350000
//PLOTLOG DD DUMMY 01360000
// 01370000

```

JCL file [2] .OPJCL.CNTL(GDI1CDEN)

CITDENS, 'iterative condensation' & 'basic fuel cycle'

```

//POCOB17C JOB {},CITDENS,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010081
// CLASS=C,TIME=0015 00020081
//***** 00030006
//CITATION EXEC PGM=CITATION,COND=(4,LT) 00040000
/**PSYS DD SUBSYS=(VPCS) 00050027
//STEPLIB DD DSN=POC3A12.CITDENS.LOAD,DISP=SHR,LABEL=(, , , IN) 00060000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(20,2)), 00070000
DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00080000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00090000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00100000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00110000
/**
***** INPUT CITATION DATA0120010
//FT05F001 DD DSN=POCOB17.PQINT.DATA(CIT),DISP=SHR,LABEL=(, , , IN) 00130057
----- 00140010
//FT06F001 DD SYSOUT=* 00150000
//FT07F001 DD UNIT=SYSDA,SPACE=(TRK,(50,50)),DCB=*.FT01F001 00160000
/**
***** INPUT 7 GROUP MICRO DATA00170010
//FT08F001 DD DSN=POCOB17.OBHZHBC3.MIC18G3,DISP=SHR,LABEL=(, , , IN) 00180088
----- 00190010
/**
***** ( OUTPUT FLUX MAP )-----**
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00260000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00270000
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00280000
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00290000
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00300000
/**
***** ( RESTART FILE )-----**
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00310000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00320000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00330000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00340000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00350000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00360000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00370000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00380000
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00390000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00400000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00410000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00420000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00430000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00440000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00450000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00460000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00470000
/**
***** ( MACRO CROSS SECTIONS )-----
//FT31F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00480000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00490000
//FT33F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00500000
/**
***** ( OUTPUT RESULT FILE )-----
//FT34F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00510000
//FT35F001 DD DUMMY 00520000
//FT36F001 DD DUMMY 00530000
//FT37F001 DD DUMMY 00540000
//FT38F001 DD DUMMY 00550000
//FT39F001 DD DUMMY 00560000
//FT40F001 DD DUMMY 00570000

```

```
//FT51F001 DD DUMMY                                00580000
/**
//FT52F001 DD DSN=POCOB17. @GDI1. DENS, DISP=(NEW, CATLG), 00591090
/**
//          SPACE=(TRK, (90, 10), RLSE), UNIT=DASD,      00592011
//          DCB=(RECFM=FBA, LRECL=137, BLKSIZE=3425)      00600065
//          *---( BURNUP FACTOR FILE )-----
//FT63F001 DD DUMMY                                00620000
//FT80F001 DD DUMMY                                00630000
//FT81F001 DD DUMMY                                00640000
//FT92F001 DD DSN=&&POFU, DISP=(NEW, PASS), SPACE=(TRK, (100, 10), RLSE), 00650000
//          UNIT=WORK, DCB=(RECFM=VSB, LRECL=4092, BLKSIZE=4096) 00660000
//          00670000
//          00680000
```

JCL file [3] .TSS.CLIST

RZOUT3 enabling MACRO

```
00000001/* CLIST CONVERSION COMPLETED 91/11/13 */
00000100PROC 0 OFF
00000201WRITE *** TSS COMMAND SYSTEM u2 Fuuu ***
00000210CONTROL NOMSG
00000300FREE F(SYSPROC)
00000400/* PROFILE PAUSE
00000500ALLOC DA('POCOB17.TSSMAC.CLIST' +
00000501 'POCOAA1.TSSMAC.CLIST' +
00000502 'POC3AA3.CMDPROC.CLIST' +
00000503 'POC3C##.CMDPROC.CLIST' +
00000610 'SYS9.CMDPROC.CLIST') +
00000700 F(SYSPROC) SHR
00000710 IF &LASTCC ^=0 THEN +
00000720 DO
00000730 WRITE *** TSS COMMAND u H JuGu u u u3 ***
00000731 SE '***PLEASE FREE YOUR DATASET *** FROM' U(POC3CA1) L
00000740 END
00000800CONTROL MSG
00000801/*EX 'POC3AA3.CMDPROC.CLIST(TSSOFF)' 'ON'
00000900EXIT
```

JCL file [4] .TSSMAC.CLIST(RZOUT3)

RZOUT3 command MACRO

```
00010045 PROC 1 MEM STM(&SYSTIME) BNK(' ')
00020053 CONTROL NOLIST MSG NOF
00030000 SET &DCB1=&STR(@1&SUBSTR(1:2,&STM)&SUBSTR(4:5,&STM)&SUBSTR(7:8,&STM))
00040000 SET &DCB2=&STR(@2&SUBSTR(1:2,&STM)&SUBSTR(4:5,&STM)&SUBSTR(7:8,&STM))
00050000 SET &DCB3=&STR(@3&SUBSTR(1:2,&STM)&SUBSTR(4:5,&STM)&SUBSTR(7:8,&STM))
00060000 SET &WORK=&STR(@W&SUBSTR(1:2,&STM)&SUBSTR(4:5,&STM)&SUBSTR(7:8,&STM))
00070000 ATTR &DCB1 REC(F B) LR( 80) BLK(3120)
```

```
00080000 ATTR &DCB2 REC(F B A) LR(133) BLK(3990)
00090000 ATTR &DCB3 INPUT
00100000 SET &SPC = &STR(&BNK&BNK&BNK&BNK&BNK&BNK&BNK&BNK)
00110048 SET &DNS = &STR(POCOB17.@MEM..DENS)
00120000 SET &LEN = &LENGTH(&DNS)
00121044 SET &MEMA = &STR(&MEM.RZ)
00140000 WRITE DNS = &DNS
00141046 WRITE MEMA = &MEMA
00150053 CONTROL NOLIST NOMSG NOF
00160098 EDIT 'POCOB17.OPEDID.DATA(&MEMA)' DATA
00170016 TOP;F 'ISTEP';L *
00180016 END NOSAVE
00190067 CONTROL LIST MSG NOF
00210067 ALLOC F(FT60F001) DA(&WORK) SP(10) T UNI(DASD) REU US(&DCB1)
00220000 OPENFILE FT60F001 OUTPUT
00230000 SET &FT60F001 = &STR(&DNS&SUBSTR(&LEN+1:80,&SPC))
00240000 PUTFILE FT60F001
00250000 CLOSFILE FT60F001
00260039 ALLOC F(FT01F001) DA('&DNS') SHR REU US(&DCB3)
00270098 ALLOC F(FT05F001) DA('POCOB17.OPEDID.DATA(&MEM.RZ)') SHR REU US(&DCB3)
00280006 ALLOC F(FT06F001) DA(@@.LIST) NE CA SP(1 1) T REU US(&DCB2)
00290000 ALLOC F(FT60F001) DA(&WORK) MOD REU
00320002 FORT 'POCOAA1.UTY.FORT77(RZOUT3NN)'
00330000 FREE F(FT01F001,FT60F001)
00340000 ALC 5 * ; ALC 6 *
00360019 WRITENR ENTER MEMBER NAME : &MEM
00370019 READ &MEMB
00380019 IF &MEMB = &STR() THEN SET &MEMB = &MEM
00381044 SET &MEMB = &STR(&MEMB.RZ)
00390099 COPY &WORK OFFPORT.DATA(&MEMB)
00400000 DEL &WORK
00410012 DEL @@.LIST
00420000 EXIT
```

JCL file [5] .OPJCL.CNTL(GDG0SJC)

SLAROM-JOINT-CITATION, 'iterative condensation'

```
//POCOB17Z JOB (), 'NEMC7GOR', MSGCLASS=X, NOTIFY=POCOB17, MSGLEVEL=(2, 0), 00010040
// CLASS=C, TIME=0010 00020033
//***** 00030000
//*** SLAROM *** 00040000
//***** 00050000
//DATAP1 EXEC PGM=DATAP1 00060019
//STEPLIB DD DSN=POC3AA3.LIB.LOAD, DISP=SHR, LABEL=(, , IN) 00070000
//FT06F001 DD SYSOUT=* 00080020
//          ***** INPUT SLAROM N.D. 00110002
//FP05F001 DD DSN=POCOB17.OPDEND.DATA(GDG0S), DISP=SHR, 00120050
//          ***** 00121002
//          LABEL=(, , IN) 00130000
//***** 00140000
//SLAGO EXEC PGM=SLAROM 00160000
//STEPLIB DD DSN=POCOEA1.SLAROM.NEDAC.LOAD, DISP=SHR 00180001
//*----- ( INPUT PDS )----- 00190000
```

```

//PDSIN DD DUMMY 00200000
/*----- ( OUTPUT PDS )----- 00210000
//PDSOUT DD DSN=&&XSG70,UNIT=WORK, 00220000
// DISP=(NEW,PASS), 00230000
// DCB=(RECFM=U,BLKSIZE=26793), 00240000
// SPACE=(CYL,(20,10,40),RLSE) 00250000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)), 00270000
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096) 00280000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00290000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00300000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00310000
//* ***** INPUT SLAROM N.D.00320002
//FT05F001 DD DSN=POCOB17.OPEND.DATA(GDGOS),DISP=SHR, 00330050
//* ----- 00331002
// LABEL=(, , IN) 00340000
//FT06F001 DD DSN=&&LIST2,DISP=(NEW,PASS), 00351031
// SPACE=(CYL,(10,3)), 00360031
// DCB=(RECFM=PBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 00370031
//FT07F001 DD DUMMY 00380000
//* ***** INPUT REF. 70 GROUP MICRO00390002
//FT08F001 DD DSN=POCOH##.JFS3J3.Y9406,DISP=SHR,LABEL=(, , IN) 00420010
//* ----- 00430002
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00440000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00450000
//FT11F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00460000
//FT12F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00470000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00480000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00490000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00500000
//FT22F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00510000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00520000
//FT26F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00530000
//FT30F001 DD DUMMY 00540000
//FT40F001 DD SYSOUT=* 00550000
//FT41F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00560000
//FT42F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00570000
//FT43F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00580000
//FT55F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00590000
//FT56F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00600000
//FT59F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00610000
//FT60F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00620000
//FT70F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00630000
//FT72F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00640000
//FT73F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00650000
//FT71F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00660000
//FT80F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00670000
//FT90F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00680000
//FT96F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00690000
//FT99F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00700000
//***** 00710000
//LOUT2 EXEC COMPACT,PARM='TYPE2,CC=NO' 00720031
//UTYIN DD DSN=&&LIST2,DISP=(OLD,DELETE) 00730031
//***** 00770000
//* JOINT ---> CITATION *//00790000
//***** 00810000
//DATAP2 EXEC PGM=DATAP1 00820019
//STEPLIB DD DSN=POC3AA3.LIB.LOAD,DISP=SHR,LABEL=(, , IN) 00830000
//* ***** INPUT CIT. DATA 70 GP.00840002
//FT05F001 DD DSN=POCOB17.OPCITD.DATA(PHAIN70), 00841034

```



```

24 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31
32
0 0
0 0
235 288 0 6 0 29 0
600MW/60CM HQ PU 40.25/45.03%,165%FUEL,4*9.75M.,24/8P ZRH/B4C OPT.
/*
//FT20F001 DD DUMMY RES CAN
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
//FT07F001 DD SYSOUT=*,FLASH=PLK1,DCB=(RECFM=UA,BLKSIZE=900,OPTCD=U)
//FT09F001 DD SYSOUT=*,FLASH=PLK2,DCB=(RECFM=UA,BLKSIZE=900,OPTCD=U)
//FT10F001 DD SYSOUT=*,FLASH=PLK3,DCB=(RECFM=UA,BLKSIZE=900,OPTCD=U)
//FT11F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
/*
*****
INPUT MULTIPLE N.D.FILE
//FT08F001 DD DSN=POCOB17.GDI1.DENS,DISP=SHR,LABEL=(, , , IN)
/*
//FT98F001 DD DSN=POCOAA1.ORGBASE.DATA,DISP=SHR,LABEL=(, , , IN)
//FT99F001 DD DUMMY
//

```

JCL file [7] .OPJCL.CNTL(GDI1SC)

SLAROM-CITATION, 'basic fuel cycle'

```

//POCOB17D JOB (),'NEMC7GOR',MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010005
// CLASS=C,TIME=0015 00020010
/****** 00030000
**** SLAROM *** SOC RODS 0% HET. RODS 00040016
***** 00050000
//DATAP1 EXEC PGM=DATAP1 00060000
//STEPLIB DD DSN=POC3AA3.LIB.LOAD,DISP=SHR,LABEL=(, , , IN) 00070000
//FT06F001 DD SYSOUT=* 00080000
/* ***** INPUT SLAROM N.D. 00110000
//FT05F001 DD DSN=POCOB17.OPDEND.DATA(GDI1S),DISP=SHR, 00120018
/* 00121000
LABEL=(, , , IN) 00130000
***** 00140000
//SLAGO EXEC PGM=SLAROM 00160000
//STEPLIB DD DSN=POCOEAL.SLAROM.NEDAC.LOAD,DISP=SHR 00180000
/*----- ( INPUT PDS )----- 00190000
//PDSIN DD DUMMY 00200000
/*----- ( OUTPUT PDS )----- 00210000
//PDSOUT DD DSN=&&XSG70,UNIT=WORK, 00220000
// DISP=(NEW,PASS), 00230000
// DCB=(RECFM=U,BLKSIZE=26793), 00240000
// SPACE=(CYL,(20,10,40),RLSE) 00250000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)), 00270000
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096) 00280000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00290000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00300000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00310000
/* ***** INPUT SLAROM N.D. 00320000
//FT05F001 DD DSN=POCOB17.OPDEND.DATA(GDI1S),DISP=SHR, 00330018
/* 00331000

```

```

// LABEL=(, , , IN) 00340000
//FT06F001 DD DSN=&&LIST2,DISP=(NEW,PASS), 00351007
// SPACE=(CYL,(10,3)), 00360007
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 00370007
//FT07F001 DD DUMMY 00380000
/* ***** INPUT REF. 70 GROUP MICRO 00390000
//FT08F001 DD DSN=POCOH#.JFS3J3.Y9406,DISP=SHR,LABEL=(, , , IN) 00420000
/* ----- 00430000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00440000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00450000
//FT11F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00460000
//FT12F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00470000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00480000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00490000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00500000
//FT22F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00510000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00520000
//FT26F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00530000
//FT30F001 DD DUMMY 00540000
//FT40F001 DD SYSOUT=* 00550000
//FT41F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00560000
//FT42F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00570000
//FT43F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00580000
//FT55F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00590000
//FT56F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00600000
//FT59F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00610000
//FT60F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00620000
//FT70F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00630000
//FT72F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00640000
//FT73F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00650000
//FT71F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00660000
//FT80F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00670000
//FT90F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00680000
//FT96F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00690000
//FT99F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00700000
/****** 00710000
***** SUPPRESS OUTPUT ***** 00711009
/*LOUT2 EXEC COMPACT,PARM='TYPE2,CC=NO' 00720009
/*UTYIN DD DSN=&&LIST2,DISP=(OLD,DELETE) 00730009
/****** 00770000
/* JOINT ---> CITATION *//00790000
/****** 00810000
//DATAP2 EXEC PGM=DATAP1 00820000
//STEPLIB DD DSN=POC3AA3.LIB.LOAD,DISP=SHR,LABEL=(, , , IN) 00830000
/* ***** INPUT CIT. DATA 70 GP. 00840000
//FT05F001 DD DSN=POCOB17.OPCITD.DATA(PEE1D700), 00841000
// 00842000
// DISP=SHR,LABEL=(, , , IN) 00850000
//FT06F001 DD SYSOUT=* 00860000
/* 00890000
/* JOINT UPDATE PROCESS *//00900001
/* 00920000
//FORT EXEC FORTECL,PARM.FORT='NOSOURCE,LC(99)' 00930001
//FORT.SYSIN DD * 00940001
DIMENSION A(500000) 00941001
LIMIT = 500000 00942001
CALL MAIN1 (A,LIMIT) 00943001
STOP 00944001
END 00945001

```



```

//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(2,5)),DCB=*.FT01F001 01440000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01450000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01460000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01470000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01480000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01490000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01500000
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01510000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01520000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01530000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01540000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01550000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01560000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01570000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01580000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01590000
/*---( MACRO CROSS SECTIONS )-----01600000
//FT31F001 DD DSN=*.JOINT1.FT09F001,DISP=(OLD,DELETE) 01610000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01620000
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01630000
/*
*****
//FT34F001 DD DSN=POCOB17.GDI3P.FT34G70N,DISP=(NEW,CATLG), 01640019
/*
UNIT=DASD,SPACE=(TRK,(10,10),RLSE),DCB=*.FT01F001 01651019
//FT51F001 DD SYSOUT=* 01652029
/*T51F001 DD DSN=&&LIST5,DISP=(NEW,PASS), 01660031
SPACE=(CYL,(10,3)), 01661031
DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 01670031
/*
01680031
/*OUT5 EXEC COMPACT,PARM='TYPE2,CC=YES' 01690000
/*TYIN DD DSN=&&LIST5,DISP=(OLD,DELETE) 01700031
/*
( JOINT.MIC ) 01710031
//JOINT2 EXEC PGM=MAIN 01720000
//STEPLIB DD DSN=&&GOSET,DISP=(OLD,PASS),LABEL=(, , IN) 01730028
//SYSPRINT DD SYSOUT=* 01740028
/*
*****
//USERPDS DD DSN=POCOB17.GDI3P70.FLUXN,DISP=(NEW,CATLG),UNIT=DASD, 01770000
/*
DCB=(RECFM=FB,LRECL=4,BLKSIZE=4000), 01780002
SPACE=(TRK,(9,3,9),RLSE) 01790000
//FT01F001 DD DSN=&&CIT,DISP=(OLD,DELETE) 01800029
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 01810000
/*
*****
//FT05F001 DD DSN=POCOB17.OPEDID.DATA(PIJ3PF),DISP=SHR 01820000
/*
*****
//FT06F001 DD SYSOUT=* 01830002
//FT90S001 DD SYSOUT=* 01840000
/*
*****
//JOINT3 EXEC PGM=MAIN 01850000
//STEPLIB DD DSN=&&GOSET,DISP=(OLD,DELETE),LABEL=(, , IN) 01870000
//SYSPRINT DD SYSOUT=* 01880028
/*
*****
//PDSIN DD DSN=POCOB17.GDI3P.PDS70G,DISP=SHR 01890028
/*
*****
//PDSOUT DD DSN=POCOB17.GDI3P18.PDSMICN, 01920000
/*
*****
UNIT=DASD,DCB=(RECFM=FB,LRECL=4,BLKSIZE=4000),DISP=(NEW,CATLG), 01921015
SPACE=(TRK,(9,3,60)) 01930032
01931015
01940018
01960032
01961002
01970000
01980029

```

```

*****
//USERPDS DD DSN=POCOB17.GDI3P70.FLUXN,DISP=SHR 02000032
/*
*****
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 02001002
/*
*****
//FT05F001 DD DSN=POCOB17.OPEDID.DATA(PIJ3PC),DISP=SHR, 02010000
/*
*****
LABEL=(, , IN) 02020002
//FT06F001 DD SYSOUT=* 02030000
//FT90S001 DD SYSOUT=* 02040000
02050000
02410000

```

JCL file [10] .OPJCL.CNTL(GDI3P70V)

CITATION, void, 70 group, 'perturbation'

```

//POCOB17P JOB ( ),JCTFBR,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010053
/*
CLASS=C,TIME=0025 00020055
/*
00021056
/*
JOINT UPDATE PROCESS *//00022056
/*
00023056
//FORT EXEC FORTECL,PARM.FORT='NOSOURCE,LC(99)' 00024056
//FORT.SYSIN DD * 00025056
DIMENSION A(500000) 00026056
LIMIT = 500000 00027056
CALL MAIN1 (A,LIMIT) 00028056
STOP 00029056
END 00029156
00029256
/*
*****
//STEPLIB DD DSN=&&GOSET,DISP=(OLD,PASS),LABEL=(, , IN) 00029356
//LKED.OLDLDM DD DSN=POCOH#.JOINT.LOAD,DISP=SHR,LABEL=(, , IN) 00029456
//LKD.SYSIN DD * 00029556
INCLUDE OLDLDM(JOINT) 00029656
ENTRY MAIN 00029756
/*
00029856
//JOINT1 EXEC PGM=MAIN 00029956
//STEPLIB DD DSN=&&GOSET,DISP=(OLD,PASS),LABEL=(, , IN) 00030056
/*
*****
//JOINT ---> CITATION *//00040000
*****
//SYSPRINT DD SYSOUT=* //00050000
/*
*****
//USERPDS DD DSN=POCOB17.GDI3P.PDS70G,DISP=SHR 00100000
18 GP. PDS (MIC.&MAC.) 00101046
//FT04F001 DD DISP=(,PASS),UNIT=WORK,SPACE=(TRK,(100,10)), 00110059
DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 00111046
/*
*****
//FT05F001 DD DSN=POCOB17.OPCITD.DATA(PIJ3PV70),DISP=SHR,LABEL=(, , IN) 00120000
18 GP. CIT. INPUT DATA 00130000
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00131046
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00140058
SPACE=(TRK,(10,10)),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 00141046
//FT09F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00150000
SPACE=(TRK,(10,10)) 00180000
00200000
00210000

```

```

//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00220000
//          SPACE=(TRK,(10,10)), 00230000
//          DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 00240000
//FT20F001 DD DUMMY 00250000
//FT50F001 DD SYSOUT=* 00260000
//*----- ( COMPACT )----- 00270000
//*OUT4 EXEC COMPACT,PARM='TYPE2,CC=NO' 00280000
//*TYIN DD DSN=&&LIST4,DISP=(OLD,DELETE) 00290000
//*----- ( CITATION )----- 00300000
//CITATION EXEC PGM=CITATION,COND=(4,LT) 00310000
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR,LABEL=(, , , IN) 00320000
//SYSPRINT DD SYSOUT=* 00330000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 00340000
//          DCB=(RECFM=VSB,LRECL=32000,BLKSIZE=32004) 00350000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00360000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00370000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00380000
//FT05F001 DD DSN=*.JOINT1.FT08F001,DISP=(OLD,DELETE) 00390000
//FT06F001 DD SYSOUT=* 00400000
//*--- ( OUTPUT ZONE AVERAGED FLUX & AXIAL BUCKLING )-----** 00410000
//FT07F001 DD DUMMY 00420000
//FT08F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 00430000
//          DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00440000
//*--- ( OUTPUT FLUX MAP )-----** 00450000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00460000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00470000
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00480000
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00490000
//*--- ( RESTART FILE )-----** 00500000
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00510000
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00520000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(2,5)),DCB=*.FT01F001 00530000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00540000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00550000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00560000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00570000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00580000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00590000
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00600000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00610000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00620000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00630000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00640000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00650000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00660000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00670000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00680000
//*--- ( MACRO CROSS SECTIONS )----- 00690000
//FT31F001 DD DSN=*.JOINT1.FT09F001,DISP=(OLD,DELETE) 00700000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00710000
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00720000
//*          ++++++*****          OUTPUT 18 GP. FLUX 00730046
//FT34F001 DD DSN=POCOB17.@GDI3P.FT34G70V,DISP=(NEW,CATLG), 00740059
//*          00741046
//          UNIT=DASD,SPACE=(TRK,(20,10),RLSE),DCB=*.FT01F001 00750057
//FT51F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00760000
//          00770000

```

JCL file [11] .OPJCL.CNTL(GDI3P18N)

CITATION, normal, 18 groups, 'perturbation'

```

//POCOB17P JOB ( ),JCTFBR,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010051
//          CLASS=C,TIME=0025 00020052
//*          JOINT UPDATE PROCESS *//00021053
//*          00023053
//FORT EXEC FORTECL,PARM.FORT='NOSOURCE,LC(99)' 00024053
//FORT.SYSIN DD * 00025053
//          DIMENSION A(500000) 00026053
//          LIMIT = 500000 00027053
//          CALL MAIN1 (A,LIMIT) 00028053
//          STOP 00029053
//          END 00029153
//*          00029253
//*--- ( NEW LOAD MODULE )----- 00029353
//LKED.OLDLM DD DSN=POCOH##.JOINT.LOAD,DISP=SHR,LABEL=(, , , IN) 00029453
//LKED.SYSIN DD * 00029553
//          INCLUDE OLDDLM(JOINT) 00029653
//          ENTRY MAIN 00029753
//*          00029853
//JOINT1 EXEC PGM=MAIN 00029953
//STEPLIB DD DSN=&&GOSET,DISP=(OLD,DELETE),LABEL=(, , , IN) 00030053
//***** 00030153
//*          JOINT ---> CITATION *//00040000
//SYSPRINT DD SYSOUT=* 00050000
//*          ++++++*****          INPUT 18 GP. MIC&MAC(PDS) 00101045
//USERPDS DD DSN=POCOB17.@GDI3P18.PDSMICN,DISP=SHR 00110057
//*          00111045
//FT04F001 DD DISP=(,PASS),UNIT=WORK,SPACE=(TRK,(100,10)), 00120000
//          DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 00130000
//*          ++++++*****          INPUT 18 GP. CIT. DATA 00131045
//FT05F001 DD DSN=POCOB17.OPCITD.DATA(PIJ3PADJ),DISP=SHR,LABEL=(, , , IN) 00140056
//*          00141045
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00150000
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00180000
//          SPACE=(TRK,(10,10)),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 00190000
//FT09F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00200000
//          SPACE=(TRK,(10,10)) 00210000
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 00220000
//          SPACE=(TRK,(10,10)), 00230000
//          DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 00240000
//FT20F001 DD DUMMY 00250000
//FT50F001 DD SYSOUT=* 00260000
//*----- ( COMPACT )----- 00270000
//*OUT4 EXEC COMPACT,PARM='TYPE2,CC=NO' 00280000
//*TYIN DD DSN=&&LIST4,DISP=(OLD,DELETE) 00290000
//*----- ( CITATION )----- 00300000
//CITATION EXEC PGM=CITATION,COND=(4,LT) 00310000
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR,LABEL=(, , , IN) 00320000
//SYSPRINT DD SYSOUT=* 00330000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 00340000
//          DCB=(RECFM=VSB,LRECL=32000,BLKSIZE=32004) 00350000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00360000

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```

//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00370000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00380000
//FT05F001 DD DSN=*.JOINT1.FT08F001,DISP=(OLD,DELETE) 00390000
//FT06F001 DD SYSOUT=* 00400000
/*----( OUTPUT ZONE AVERAGED FLUX & AXIAL BUCKLING )-----** 00410000
//FT07F001 DD DUMMY 00420000
//FT08F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 00430000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 00440000
/*----( OUTPUT FLUX MAP )-----** 00450000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00460000
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00470000
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00480000
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00490000
/*----( RESTART FILE )-----** 00500000
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00510000
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00520000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(2,5)),DCB=*.FT01F001 00530000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00540000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00550000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 00560000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 00570000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00580000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00590000
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00600000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00610000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00620000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00630000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00640000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00650000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00660000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00670000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 00680000
/*----( MACRO CROSS SECTIONS )-----** 00690000
//FT31F001 DD DSN=*.JOINT1.FT09F001,DISP=(OLD,DELETE) 00700000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00710000
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 00720000
/* ***** OUTPUT 18 GP. FLUX 00730045
//FT34F001 DD DSN=POCOB17. @GDI3P.FT34G18N,DISP=(NEW,CATLG), 00740057
/* 00741045
// UNIT=DASD,SPACE=(TRK,(10,10),RLSE),DCB=*.FT01F001 00750054
//FT51F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00760000
// 00770000

```

```

DIMENSION A(500000) 00080027
LIMIT = 500000 00090027
CALL MAIN1 (A,LIMIT) 00100027
STOP 00110027
END 00120027
/* 00130027
/*----( NEW LOAD MODULE )-----** 00140027
//LKED.OLDLIM DD DSN=POCOH##.JOINT.LOAD,DISP=SHR,LABEL=(,,IN) 00150027
//LKED.SYSIN DD * 00160027
INCLUDE OLDLIM(JOINT) 00170027
ENTRY MAIN 00180027
/* 00190027
//JOINT1 EXEC PGM=MAIN 00200027
//STEPLIB DD DSN=*&GOSET,DISP=(OLD,PASS),LABEL=(,,IN) 00210027
/****** 00770000
/* JOINT --> CITATION * 00790000
/****** 00810000
//SYSPRINT DD SYSOUT=* 00970000
/* ***** INPUT 70 GP. MIC&MAC 00971017
//USERPDS DD DSN=POCOB17. @GDI3P.PDS70G,DISP=SHR 00980030
/* 00981017
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,10)), 00990000
// DISP=(,PASS),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 01000000
/* ***** INPUT CIT. DATA 70 GP. 01010002
//FT05F001 DD DSN=POCOB17.OPCITD.DATA(PIJ3P70D),DISP=SHR, 01011029
/* 01012002
// LABEL=(,,IN) 01020000
//FT06F001 DD DSN=*&LIST4,DISP=(NEW,PASS), 01030000
// SPACE=(CYL,(10,3)), 01040000
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 01050000
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01060000
// SPACE=(TRK,(10,10)), 01070000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01080000
//FT09F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01090000
// SPACE=(TRK,(10,10)) 01100000
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01110000
// SPACE=(TRK,(10,10)), 01120000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01130000
//FT20F001 DD DUMMY 01140000
//FT50F001 DD SYSOUT=* 01150000
/* 01160000
//LOUT4 EXEC COMPACT,PARM='TYPE2,CC=NO' 01170000
//UTYIN DD DSN=*&LIST4,DISP=(OLD,DELETE) 01180000
/*----( CITATION )-----** 01190000
//CITATION EXEC PGM=CITATION,COND=(4,LT) 01200000
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR 01210000
//SYSPRINT DD SYSOUT=* 01220000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01230000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 01240000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01250000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01260000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01270000
//FT05F001 DD DSN=*.JOINT1.FT08F001,DISP=(OLD,DELETE) 01280000
//FT06F001 DD SYSOUT=* 01290000
/*----( OUTPUT ZONE AVERAGED FLUX & AXIAL BUCKLING )-----** 01300000
//FT07F001 DD DSN=*&CIT,DISP=(NEW,PASS),UNIT=WORK, 01310000
// SPACE=(TRK,(2,1),RLSE), 01320000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200) 01330000
//FT08F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01340000

```

JCL file [12] .OPJCL.CNTL(GDI3P70D)

CITATION, Doppler, 70 groups, 'perturbation'

```

//POCOB17P JOB (),'NEMC7GOR',MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010025
// CLASS=C,TIME=0025 00020026
/* 00030027
/* JOINT UPDATE PROCESS **/00040027
/* 00050027
//FORT EXEC FORTECL,PARM.FORT='NOSOURCE,LC(99)' 00060027
//FORT.SYSIN DD * 00070027

```



```

//FT40F001 DD DISP=(NEW,PASS,DELETE),          *DELAYED NEUTRON * 00320000
//          SPACE=(TRK,(10,5)),UNIT=WORK,        *FOR NEXT JOB STEP* 00330000
//          DSN=&&BDELAY                            *BY BINARY FORM * 00340000
//*****
//FT50F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137, 00350000
//          BLKSIZE=19043)                          00360000
//-----
//*          EXPAND ABLE MEMORY SIZE                00370000
//-----
// EXEC FORTECL, PARM.FORT='OPT(B),NOSTATIS,ELM(*),NOAE', 00381139
//          PARM.LKED='MAP,LIST,LET,SIZE=(512K,72K)',      00381239
//          REGION.LKED=1024K,COND.LKED=(8,LT)             00381339
//          DD DSN=POCOB17.OPFORT.FORT(PERK),DISP=SHR,LABEL=(,,IN) 00381439
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381539
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381639
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381739
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381840
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381939
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382039
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382139
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382239
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382339
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382439
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382539
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382639
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382739
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382839
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00382939
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 003900039
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00400039
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00410039
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00420000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00430015
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00440041
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00450015
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00460035
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00470000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00480000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00490000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00500000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00510000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00520000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00530000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00540000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00550000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00560000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00570000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00580000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00590000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00600000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00610000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00620000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00630019
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00640041
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00641015
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00650000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00660000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00670000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00680000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00690000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00700000
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00710011

```

```

//*****
//LSTPKY EXEC PGM=LSTPKY,COND=(4,LT)
//STEPLIB DD DSN=POCOA1.INTERF.LOAD,DISP=SHR
//*PSYS DD SUBSYS=(VPCS)
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
//FT05F001 DD DUMMY
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
//

```

```

00720011
00730009
00740009
00750009
00760009
00770009
00780009
00790000

```

JCL file [14] .OPJCL.CNTL(GDI3PDNF)

PERKY, delayed neutron fraction, 'perturbation'

```

//POCOB17P JOB ( ), PERKY,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010039
//          CLASS=C,TIME=0005                                     00020041
//* 18GP          DELAYED NEUTRON FRACTION                       00021027
//JOINT EXEC PGM=JOINT                                          00030015
//STEPLIB DD DSN=POCOA1.JOINTL.LOAD,DISP=SHR,LABEL=(,,IN)     00040025
//*PSYS DD SUBSYS=(VPCS)                                        00050000
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00060000
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00070000
//FT04F001 DD SPACE=(TRK,(50,5)),UNIT=WORK                    00080000
//          ***** INPUT PERKY DATA                          00090015
//          DD DSN=POCOB17.OPEPID.DATA(PIJ3PDNF),DISP=SHR,LABEL=(,,IN) 00100043
//          ***** INPUT 18 GP FLUX D&A00430015              00110015
//          DD DSN=POCOB17.@GDI3P.FT34G70N,DISP=SHR,LABEL=(,,IN) 00111015
//          ***** INPUT 18 GP PDS FILES                       00120045
//          DD DSN=POCOB17.@GDI3P18.PDSMICN,DISP=SHR,LABEL=(,,IN) 00130045
//          DD DSN=POCOB17.@GDI3P18.PDSMICN,DISP=SHR,LABEL=(,,IN) 00141015
//          *****
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK,                *INPUT DATA * 00150000
//          SPACE=(TRK,(50,5)),DSN=&&INPUT,                      *FOR NEXT JOB STEP* 00160000
//          DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240)                *BY CARD FORM * 00170000
//          *****
//FT09F001 DD DISP=(NEW,PASS,DELETE),                           *CROSS SECTIONS * 00200000
//          SPACE=(TRK,(50,5)),UNIT=WORK,                       *FOR NEXT JOB STEP* 00210000
//          DSN=&&BMACRO                                           *BY BINARY FORM * 00220000
//          *****
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK,                *CROSS SECTIONS * 00240000
//          SPACE=(TRK,(50,5)),DSN=&&CCROSS,                     *FOR NEXT JOB STEP* 00250000
//          DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240)                *BY CARD FORM * 00260000
//          *****
//FT30F001 DD DISP=(NEW,PASS,DELETE),                           *CROSS SECTIONS * 00270000
//          SPACE=(TRK,(50,5)),UNIT=WORK,                       *FOR NEXT JOB STEP* 00280000
//          DSN=&&EMICRO                                           *BY BINARY FORM * 00290000
//          *****
//FT40F001 DD DISP=(NEW,PASS,DELETE),                           *DELAYED NEUTRON * 00320000
//          SPACE=(TRK,(10,5)),UNIT=WORK,                       *FOR NEXT JOB STEP* 00330000
//          DSN=&&BDELAY                                           *BY BINARY FORM * 00340000
//          *****
//FT50F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00360000
//-----
//*          EXPAND PERKY MEMORY SIZE                            00381144
//-----
//          DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN) 00381244

```

```

-----
// EXEC FORTECL, PARM.FORT='OPT(B),NOSTATIS,ELM(*)',NOAE',
// PARM.LKED='MAP,LIST,LET,SIZE=(512K,72K)',
// REGION.LKED=1024K,COND.LKED=(8,LT)
// FORT.SYSIN DD DSN=POCOB17.OPFORT.FORT(PERK),DISP=SHR,LABEL=(,,IN)
// LKED.OLDL DD DSN=POCOH##.PERKY.LOAD,DISP=SHR,LABEL=(,,IN)
// LKED.SYSIN DD *
// INCLUDE OLDL(Y9607)
// ENTRY MAIN
// NAME MAIN
/*
-----
// GO PERKY
-----
// PERKY EXEC PGM=MAIN
// STEPLIB DD DSN=&&GOSET,DISP=(OLD,DELETE)
// *ERKY EXEC PGM=Y9308,COND=(4,LT)
// *TEPLIB DD DSN=POCOB15.PERKY.LOAD,DISP=SHR
// *PSYS DD SUBSYS=(VPCS)
// SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// ***** INPUT 18 GP FLUX D&A00430015
// FT01F001 DD DSN=POCOB17.@GDI3P.FT34G18N,DISP=SHR,LABEL=(,,IN)
// *****
// FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=(BLKSIZE=18628)
// FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT05F001 DD DSN=*.JOINT.FT08F001,DISP=(OLD,DELETE)
// FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// FT08F001 DD DSN=*.JOINT.FT09F001,DISP=(OLD,DELETE)
// FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT11F001 DD DSN=*.JOINT.FT30F001,DISP=(OLD,DELETE)
// FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT17F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT21F001 DD DSN=*.JOINT.FT40F001,DISP=(OLD,DELETE)
// FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// ***** INPUT 18 GP. UNPERT. FLUX
// FT23F001 DD DSN=POCOB17.@GDI3P.FT34G18N,DISP=SHR,LABEL=(,,IN)
// *****
// FT23F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// *****
// FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// *****
// LSTPKY EXEC PGM=LSTPKY,COND=(4,LT)
// STEPLIB DD DSN=POCOAAL.INTERF.LOAD,DISP=SHR
// *PSYS DD SUBSYS=(VPCS)
// SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// FT05F001 DD DUMMY
// FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
//

```

JCL file [15] .OPJCL.CNTL(GDI3PEV)

PERKY, Na void, 'perturbation'

```

// POCOB17P JOB (),PERKY,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0),
// CLASS=C,TIME=0005
// * 70 GP. NA VOID PERTURBATION , EXACT
// JOINT EXEC PGM=JOINT
// STEPLIB DD DSN=POCOAAL.JOINTL.LOAD,DISP=SHR,LABEL=(,,IN)
// *PSYS DD SUBSYS=(VPCS)
// SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// FT04F001 DD SPACE=(TRK,(50,5)),UNIT=WORK
// FT05F001 DD DSN=POCOB17.OPEDID.DATA(PIJ3PEV),DISP=SHR,LABEL=(,,IN)
// ***** INPUT PERKY DATA
// ***** INPUT 18 GP PDS FILES
// USERPDS DD DSN=POCOB17.@GDI3P.PDS70G,DISP=SHR,LABEL=(,,IN)
// *****
// FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK, *INPUT DATA *
// SPACE=(TRK,(50,5)),DSN=&&INPUT, *FOR NEXT JOB STEP*
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240) *BY CARD FORM *
// *****
// FT09F001 DD DISP=(NEW,PASS,DELETE), *CROSS SECTIONS *
// SPACE=(TRK,(50,5)),UNIT=WORK, *FOR NEXT JOB STEP*
// DSN=&&BMACRO *BY BINARY FORM *
// *****
// FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK, *CROSS SECTIONS *
// SPACE=(TRK,(50,5)),DSN=&&CCROSS, *FOR NEXT JOB STEP*
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240) *BY CARD FORM *
// *****
// FT30F001 DD DISP=(NEW,PASS,DELETE), *CROSS SECTIONS *
// SPACE=(TRK,(50,5)),UNIT=WORK, *FOR NEXT JOB STEP*
// DSN=&&BMICRO *BY BINARY FORM *
// *****
// FT40F001 DD DISP=(NEW,PASS,DELETE), *DELAYED NEUTRON *
// SPACE=(TRK,(10,5)),UNIT=WORK, *FOR NEXT JOB STEP*
// DSN=&&BDELAY *BY BINARY FORM *
// *****
// FT50F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// *****
// PERKY EXEC PGM=Y9308,COND=(4,LT)
// STEPLIB DD DSN=POCOAAL.PERKYL.LOAD,DISP=SHR
// *PSYS DD SUBSYS=(VPCS)
// SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// ***** INPUT 18 GP FLUX D&A00430015
// FT01F001 DD DSN=POCOB17.@GDI3P.FT34G70N,DISP=SHR,LABEL=(,,IN)
// *****
// FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=(BLKSIZE=18628)
// FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT05F001 DD DSN=*.JOINT.FT08F001,DISP=(OLD,DELETE)
// FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043)
// FT08F001 DD DSN=*.JOINT.FT09F001,DISP=(OLD,DELETE)
// FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
// FT11F001 DD DSN=*.JOINT.FT30F001,DISP=(OLD,DELETE)

```



```

//FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00540000
//FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00550000
//FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00560000
//FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00570000
//FT17F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00580000
//FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00590000
//FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00600000
//FT21F001 DD DSN=*.JOINT.FT40F001,DISP=(OLD,DELETE) 00610000
//FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00620000
//*
//***** INPUT 18 GP. PERT. FLUX
//FT23F001 DD DSN=POCOB17.@GDI3P.FT34G70V,DISP=SHR,LABEL=(, , , IN) 00630015
//*
//*****
//FT23F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00640035
//*
//*****
//FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00641015
//FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00650000
//FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00660000
//FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00670000
//
//*****
//LSTPKY EXEC PGM=LSTPKY,COND=(4,LT) 00670000
//STEPLIB DD DSN=POCOA1.INTERF.LOAD,DISP=SHR 00680000
//*PSYS DD SUBSYS=(VPCS) 00690000
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00700000
//FT05F001 DD DUMMY 00710011
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00720011
//
//*****
//LSTPKY EXEC PGM=LSTPKY,COND=(4,LT) 00730009
//STEPLIB DD DSN=POCOA1.INTERF.LOAD,DISP=SHR 00740009
//*PSYS DD SUBSYS=(VPCS) 00750009
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00760009
//FT05F001 DD DUMMY 00770009
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00780009
//
//*****

```

JCL file [16] .OPJCL.CNTL(GDI3PFD0)

PERKY, Doppler (by zone), 'perturbation'

```

//POCOB17P JOB (),PERKY,MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010027
// CLASS=C,TIME=0015 00020028
//* DOPPLER PERTURBATION, I' ST ORDER, BY REGION 00021021
//JOINT EXEC PGM=JOINT 00030015
//STEPLIB DD DSN=POCOH##.JOINT.LOAD,DISP=SHR,LABEL=(, , , IN) 00040022
//*PSYS DD SUBSYS=(VPCS) 00050000
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00060000
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00070000
//FT04F001 DD SPACE=(TRK,(50,5)),UNIT=WORK 00080000
//*
//***** INPUT PERKY DATA 00090015
//FT05F001 DD DSN=POCOB17.OPEDID.DATA(PIJ3PFD0),DISP=SHR,LABEL=(, , , IN) 00100028
//*
//***** INPUT 18 GP PDS FILES 00110015
//USERPDS DD DSN=POCOB17.@GDI3P18.PDSMICN,DISP=SHR,LABEL=(, , , IN) 00120030
// DD DSN=POCOB17.@GDI3P18.PDSMICD,DISP=SHR,LABEL=(, , , IN) 00130030
//
//***** 00141015
//***** 00150000
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK, *INPUT DATA * 00160000
// SPACE=(TRK,(50,5)),DSN=*&&INPUT, *FOR NEXT JOB STEP* 00170000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240) *BY CARD FORM * 00180000
//***** 00190000
//FT09F001 DD DISP=(NEW,PASS,DELETE), *CROSS SECTIONS * 00200000

```

```

// SPACE=(TRK,(50,5)),UNIT=WORK, *FOR NEXT JOB STEP* 00210000
// DSN=*&&BMACRO *BY BINARY FORM * 00220000
//***** 00230000
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=WORK, *CROSS SECTIONS * 00240000
// SPACE=(TRK,(50,5)),DSN=*&&CCROSS, *FOR NEXT JOB STEP* 00250000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=6240) *BY CARD FORM * 00260000
//***** 00270000
//FT30F001 DD DISP=(NEW,PASS,DELETE), *CROSS SECTIONS * 00280000
// SPACE=(TRK,(50,5)),UNIT=WORK, *FOR NEXT JOB STEP* 00290000
// DSN=*&&BMICRO *BY BINARY FORM * 00300000
//***** 00310000
//FT40F001 DD DISP=(NEW,PASS,DELETE), *DELAYED NEUTRON * 00320000
// SPACE=(TRK,(10,5)),UNIT=WORK, *FOR NEXT JOB STEP* 00330000
// DSN=*&&BDELAY *BY BINARY FORM * 00340000
//***** 00350000
//FT50F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00360000
//***** 00370000
//PERKY EXEC PGM=Y9607,COND=(4,LT) 00380029
//STEPLIB DD DSN=POCOH##.PERKY.LOAD,DISP=SHR 00400029
//*PSYS DD SUBSYS=(VPCS) 00410000
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00420000
//***** INPUT 18 GP FLUX D&A00430015
//FT01F001 DD DSN=POCOB17.@GDI3P.FT34G18N,DISP=SHR,LABEL=(, , , IN) 00440030
//*
//***** 00450015
//FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=(BLKSIZE=18628) 00460000
//FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00470000
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00480000
//FT05F001 DD DSN=*.JOINT.FT08F001,DISP=(OLD,DELETE) 00490000
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00500000
//FT08F001 DD DSN=*.JOINT.FT09F001,DISP=(OLD,DELETE) 00510000
//FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00520000
//FT11F001 DD DSN=*.JOINT.FT30F001,DISP=(OLD,DELETE) 00530000
//FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00540000
//FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00550000
//FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00560000
//FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00570000
//FT17F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00580000
//FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00590000
//FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00600000
//FT21F001 DD DSN=*.JOINT.FT40F001,DISP=(OLD,DELETE) 00610000
//FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00620000
//*
//***** INPUT 18 GP. UNPERT. FLUX 00630019
//FT23F001 DD DSN=POCOB17.@GDI3P.FT34G18N,DISP=SHR,LABEL=(, , , IN) 00640030
//*
//***** 00650000
//FT23F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00650000
//***** 00660000
//FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00670000
//FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00680000
//FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00690000
//FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001 00700000
//
//***** 00710011
//LSTPKY EXEC PGM=LSTPKY,COND=(4,LT) 00720011
//STEPLIB DD DSN=POCOA1.INTERF.LOAD,DISP=SHR 00730009
//*PSYS DD SUBSYS=(VPCS) 00740009
//SYSPRINT DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00750009
//FT05F001 DD DUMMY 00760009
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 00770009
//
//***** 00790000

```



```

C   MAXM ----- TOTAL LENGTH OF MACROSCOPIC CROSS SECTION DATA 00001500
C   MAXI ----- TOTAL LENGTH OF MICROSCOPIC CROSS SECTION DATA 00001600
C   NMIC ----- NO. OF MATERIAL WHICH IS INCLUDED IN NMICRO 00001700
C   NMICRO(NMIC) ---- MATERIAL CODE NAME (OUTPUT) 00001800
C   NUCN ----- NUCLEAR NO. 00001900
C   NMNUC(NUCN) ---- MATERIAL CODE NAME (ALL) 00002000
C   DENHOM(NUCN) ---- HOMOGENIZED NO. DENSITY (ALL) 00002100
C   IDATE(2) ----- PROCESSED DATE (CHARACTOR) 00002200
C   ENBND ----- ENERGY BOUNDARY 00002300
C   00002400
C   DIMENSION NMPDS(2),NMCELL(2),ITITLE(18),NMICRO(1),NMNUC(1) 00002500
C   DIMENSION DENHOM(1),IDATE(2),MEMB(2),IWORK(250),WORK(250) 00002601
C   DIMENSION DENHOM(1),IDATE(2),MEMB(2),IWORK(500),WORK(500) 00002701
C   DIMENSION ENBND(1) 00002801
C   EQUIVALENCE (IWORK(1),WORK(1)) 00002901
C+  DATA MAXL /250/ (93/07/06) 00003001
C   DATA MAXL /500/ 00003101
C   00003201
C   LENG = 27+NMIC+2*NUCN 00003301
C   IF (LENG.GT.MAXL) GO TO 7100 00003401
C   CALL NAMSET (NMCELL,NMPROG,NMIC,NMICRO,1,MEMB) 00003501
C   00003601
C   RECORDING DATA SETUP 00003701
C   00003801
C   IWORK(1) = MAXG 00003901
C   IWORK(2) = IDS 00004001
C   IWORK(3) = IPL 00004101
C   IWORK(4) = MAXM 00004201
C   IWORK(5) = MAXI 00004301
C   IWORK(6) = NMIC 00004401
C   IWORK(7) = NUCN 00004501
C   CALL DATE (IWORK(8)) 00004601
C   DO 10 I=1,18 00004701
C   IWORK(I+9) = ITITLE(I) 00004801
10  CONTINUE 00004901
C   IF (NMIC.LE.0) GO TO 30 00005001
C   DO 20 I=1,NMIC 00005101
C   IWORK(I+27) = NMICRO(I) 00005201
20  CONTINUE 00005301
C   00005401
30  CONTINUE 00005501
C   NW = 27 + NMIC 00005601
C   IF (NUCN.LE.0) GO TO 60 00005701
C   DO 40 I=1,NUCN 00005801
C   IWORK(I+NW) = NMNUC(I) 00005901
40  CONTINUE 00006001
C   NW = NW+NUCN 00006101
C   DO 50 I=1,NUCN 00006201
C   WORK(I+NW) = DENHOM(I) 00006301
50  CONTINUE 00006401
C   NW = NW+NUCN 00006501
60  CONTINUE 00006601
C   DO 61 I=1,MAXG+1 00006701
C   NW = NW +1 00006801
C   WORK(NW) = ENEND(I) 00006901
61  CONTINUE 00007001
C   CALL PDSPUT (NMPDS,MEMB,IWORK,NW) 00007101
C   GO TO 9999 00007201
C   00007301
C   ENTRY GETIDX(NMPDS,NMCELL,NMPROG,ITITLE,MAXG,IDS,IPL,MAXM,
+   MAXI,NMIC,NMICRO,NUCN,NMNUC,DENHOM,IDATE,

```

```

+   ENEND) 00007401
C   00007501
C   READ ID DATA FROM PDS FILE 00007601
C   NMIC = 1 00007701
C   CALL NAMSET (NMCELL,NMPROG,NMIC,NMICRO,1,MEMB) 00007801
C   CALL PDSLEN (NMPDS,MEMB,LENG) 00007901
C   IF (LENG.LT.0) GO TO 7000 00008001
C   IF (LENG.GT.MAXL) GO TO 7100 00008101
C   CALL PDSGET (NMPDS,MEMB,IWORK,LENG) 00008201
C   MAXG = IWORK(1) 00008301
C   MAXGW = IWORK(1) 00008401
C   IDS = IWORK(2) 00008501
C   IPL = IWORK(3) 00008601
C   MAXM = IWORK(4) 00008701
C   MAXI = IWORK(5) 00008801
C   NMIC = IWORK(6) 00008901
C   NMICW = IWORK(6) 00009001
C   NUCN = IWORK(7) 00009101
C   NUCNW = IWORK(7) 00009201
C   IDATE(1) = IWORK(8) 00009301
C   IDATE(2) = IWORK(9) 00009401
C   DO 70 I=1,18 00009501
C   ITITLE(I) = IWORK(I+9) 00009601
70  CONTINUE 00009701
C   IF (NMICW.LE.0) GO TO 90 00009801
C   DO 80 I=1,NMICW 00009901
C   NMICRO(I) = IWORK(I+27) 00100001
80  CONTINUE 00101001
90  CONTINUE 00102001
C   IPOS = 27 + NMICW 00103001
C   IF (NUCNW.LE.0) GO TO 120 00104001
C   NW = 27+NMICW 00105001
C   DO 100 I=1,NUCNW 00106001
C   NMNUC(I) = IWORK(I+NW) 00107001
100 CONTINUE 00108001
C   NW = NW+NUCNW 00109001
C   DO 110 I=1,NUCNW 00110001
C   DENHOM(I) = WORK(I+NW) 00111001
110 CONTINUE 00112001
C   IPOS = NW + NUCNW 00113001
120 CONTINUE 00114001
C   DO 140 I=1,MAXGW+1 00115001
C   ENEND(I) = WORK(I+IPOS) 00116001
140 CONTINUE 00117001
C   GO TO 9999 00118001
C   00119001
C   ENTRY FNDIDX(NMPDS,NMCELL,NMPROG,NLENG,
*   MAXG,IDS,IPL,NMIC,NUCN) 00120001
C   00121001
C   00122001
C   FIND ID DATA FROM PDS FILE 00123001
C   MAXG = 0 00124001
C   IDS = 0 00125001
C   IPL = 0 00126001
C   NMIC = 0 00127001
C   NUCN = 0 00128001
C   CALL NAMSET (NMCELL,NMPROG,0,1,0,1,MEMB) 00129001
C   CALL PDSLEN (NMPDS,MEMB,NLENG) 00130001
C   IF (NLENG.LE.0) GO TO 9999 00131001
C   CALL PDSGET (NMPDS,MEMB,IWORK,7) 00132001

```

```

MAXG = IWORK(1)
IDS = IWORK(2)
IPL = IWORK(3)
NMIC = IWORK(6)
NUCN = IWORK(7)
GO TO 9999
C
7000 CONTINUE
WRITE (6,7010) MEMB
7010 FORMAT ('0***** ERROR MEMBER NAME ''',2A4, '' DOES NOT EXIST'')
STOP 100
C
7100 CONTINUE
WRITE (6,65) MAXL,LENG
65 FORMAT ('0***** ERROR AREA SIZE IS TOO SMALL *****',
1 /, ' ',T20,'RESERVED AREA IS ',I5,' WORDS',
2 /, ' ',T20,'REQUIRED AREA IS ',I5,' WORDS')
STOP 110
C
9999 CONTINUE
RETURN
END
SUBROUTINE PUTID (NMPDS,NMCELL,NMPROG,ITITLE,MAXG,IDS,IPL,MAXM,
+ MAXI,NMIC,NMICRO,NUCN,NMNUC,DENHOM)
C
C THIS SUBROUTINE DOES I/O OF ID DATA
C
C ARGUMENT LIST
C NMPDS(2) ----- DATA SET NAME OF PDS FILE
C NMCELL(2) ----- REGION NAME
C NMPROG ----- PROCESSED PROGRAM NAME
C ITITLE(18) ----- TITLE DATA
C MAXG ----- NO. OF ENERGY GROUP
C IDS ----- NO. OF SINK GROUP
C IPL ----- HIGHEST ORDER OF LEGENDRE EXPANSION
C MAXM ----- TOTAL LENGTH OF MACROSCOPIC CROSS SECTION DATA
C MAXI ----- TOTAL LENGTH OF MICROSCOPIC CROSS SECTION DATA
C NMIC ----- NO. OF MATERIAL WHICH IS INCLUDED IN NMICRO
C NMICRO(NMIC) ---- MATERIAL CODE NAME (OUTPUT)
C NUCN ----- NUCLEAR NO.
C NMNUC(NUCN) ---- MATERIAL CODE NAME (ALL)
C DENHOM(NUCN) ---- HOMOGENIZED NO. DENSITY (ALL)
C IDATE(2) ----- PROCESSED DATE (CHARACTOR)
C
C DIMENSION NMPDS(2),NMCELL(2),ITITLE(18),NMICRO(1),NMNUC(1)
C$ DIMENSION DENHOM(1),IDATE(2),MEMB(2),IWORK(250),WORK(250)
C DIMENSION DENHOM(1),IDATE(2),MEMB(2),IWORK(500),WORK(500)
C EQUIVALENCE (IWORK(1),WORK(1))
C+&&& DATA MAXL /250/ (93/07/06)
C DATA MAXL /500/
C
C LENG = 27+NMIC+2*NUCN
C IF (LENG.GT.MAXL) GO TO 7100
C CALL NAMSET (NMCELL,NMPROG,NMIC,NMIC,1,MEMB)
C
C RECORDING DATA SETUP
C
IWORK(1) = MAXG
IWORK(2) = IDS
IWORK(3) = IPL

```

```

00013301
00013401
00013501
00013601
00013701
00013801
00013901
00014001
00014101
00014201
00014301
00014401
00014501
00014601
00014701
00014801
00014901
00015001
00015101
00015201
00015301
00015401
00000100
00000200
00000300
00000400
00000500
00000600
00000700
00000800
00000900
00001000
00001100
00001200
00001300
00001400
00001500
00001600
00001700
00001800
00001900
00002000
00002100
00002200
00002300
00002401
00002501
00002601
00002701
00002801
00002901
00003001
00003101
00003201
00003301
00003401
00003501
00003601
00003701

```

```

IWORK(4) = MAXM
IWORK(5) = MAXI
IWORK(6) = NMIC
IWORK(7) = NUCN
CALL DATE (IWORK(8))
DO 10 I=1,18
IWORK(I+9) = ITITLE(I)
10 CONTINUE
IF (NMIC.LE.0) GO TO 30
DO 20 I=1,NMIC
IWORK(I+27) = NMICRO(I)
20 CONTINUE
30 CONTINUE
NW = 27 + NMIC
IF (NUCN.LE.0) GO TO 60
DO 40 I=1,NUCN
IWORK(I+NW) = NMNUC(I)
40 CONTINUE
NW = NW+NUCN
DO 50 I=1,NUCN
WORK(I+NW) = DENHOM(I)
50 CONTINUE
NW = NW+NUCN
60 CONTINUE
CALL PDSPUT (NMPDS,MEMB,IWORK,NW)
GO TO 9999
C
ENTRY GETID (NMPDS,NMCELL,NMPROG,ITITLE,MAXG,IDS,IPL,MAXM,
+ MAXI,NMIC,NMICRO,NUCN,NMNUC,DENHOM,IDATE)
C
C READ ID DATA FROM PDS FILE
NMIC = 1
CALL NAMSET (NMCELL,NMPROG,NMIC,NMIC,NMICRO,1,MEMB)
CALL PDSLEN (NMPDS,MEMB,LENG)
IF (LENG.LT.0) GO TO 7000
IF (LENG.GT.MAXL) GO TO 7100
CALL PDSGET (NMPDS,MEMB,IWORK,LENG)
MAXG = IWORK(1)
IDS = IWORK(2)
IPL = IWORK(3)
MAXM = IWORK(4)
MAXI = IWORK(5)
NMIC = IWORK(6)
NMICW = IWORK(6)
NUCN = IWORK(7)
NUCNW = IWORK(7)
IDATE(1) = IWORK(8)
IDATE(2) = IWORK(9)
DO 70 I=1,18
ITITLE(I) = IWORK(I+9)
70 CONTINUE
IF (NMICW.LE.0) GO TO 90
DO 80 I=1,NMICW
NMICRO(I) = IWORK(I+27)
80 CONTINUE
90 CONTINUE
IF (NUCNW.LE.0) GO TO 120
NW = 27+NMICW
DO 100 I=1,NUCNW

```

```

00003801
00003901
00004001
00004101
00004201
00004301
00004401
00004501
00004601
00004701
00004801
00004901
00005001
00005101
00005201
00005301
00005401
00005501
00005601
00005701
00005801
00005901
00006001
00006101
00006201
00006301
00006401
00006501
00006601
00006701
00006801
00006901
00007001
00007101
00007201
00007301
00007401
00007501
00007601
00007701
00007801
00007901
00008001
00008101
00008201
00008301
00008401
00008501
00008601
00008701
00008801
00008901
00009001
00009101
00009201
00009301
00009401
00009501
00009601

```

```

NMNUC(I) = IWORK(I+NW)
100 CONTINUE
NW = NW+NUCNW
DO 110 I=1,NUCNW
DENHOM(I) = WORK(I+NW)
110 CONTINUE
120 CONTINUE
GO TO 9999

C
7000 CONTINUE
WRITE (6,7010) MEMB
7010 FORMAT ('0***** ERROR MEMBER NAME ''',2A4,''' DOES NOT EXIST')
STOP 100

C
7100 CONTINUE
WRITE (6,65) MAXL,LENG
65 FORMAT ('0***** ERROR AREA SIZE IS TOO SMALL *****',
1 //, ' ',T20,'RESERVED AREA IS ',I5,' WORDS',
2 //, ' ',T20,'REQUIRED AREA IS ',I5,' WORDS')
STOP 110

C
9999 CONTINUE
RETURN
END
SUBROUTINE NAMSET (NMCELL,NMPROG,MICCOD,NMIC,NMICRO,KACC,MEMB)

C
C THIS SUBROUTINE SETS MEMBER NAME OF PDS FILE
C
C --- 83/08/11 --- ADDITION OF KACC=6 OPTION FOR ZONE AVERAGED FLUX
C --- 83/08/11 --- BY T.MITSUNARI JAIS
C
C ARGUMENT LIST
C NMCELL ---- REGION NAME (6 CHARACTER)
C NMPROG ---- PROCESSED PROGRAM NAME (4 CHARACTER)
C MICCOD ---- NUCLEAR CODE NAME (CHARACTOR OR INTEGER)
C NMIC ---- NO. OF NUCLEAR WHICH IS OUTPUTED TO PDS FILE
C NMICRO ---- NUCLEAR CODE NAME TABLE
C KACC ----- ACCESS CEDE = 1 ; ID
C = 2 ; MACROSCOPIC CROSS SECTION
C = 3 ; MICROSCOPIC CROSS SECTION
C = 4 ; BUCKLING
C = 5 ; FISSION SPECTRUM
C = 6 ; ZONE AVERAGED FLUX
C MEMB ----- MEMBER NAME OF PDS FILE
C
C DIMENSION NMCELL(2),NMICRO(NMIC)
C*-----< 1 LINE UPDATE (93/05/17)
C* CHARACTER*4 MEMB(2)
C* CHARACTER*4 MEMB(2), NUM
C*-----< 2 LINES INSERT (93/05/17)
C* INTEGER*4 SPACE
C* DATA SPACE / ' ' /
C
C INTEGER KODE(10,2)
1 / 'ESEL','TIMS','SLAR','PIGE','EXPA','ANIS','CITA',3* ' '
2 'E','T','S','P','G','A','C',3* '/'
C* DATA IDCOD, KBUCK, MACRO, KHI, LUXF
C*-----< 1 LINE UPDATE (93/05/17)
C* + /'0','@','$','#','Z' '/'

```

```

00009701
00009801
00009901
00010001
00010101
00010201
00010301
00010401
00010501
00010601
00010701
00010801
00010901
00011001
00011101
00011201
00011301
00011401
00011501
00011601
00011701
00011801
00011901
00012001
00010000
00020000
00030000
00040000
00050000
00060000
00070000
00080000
00090000
00100000
00110000
00120000
00130000
00140000
00150000
00160000
00170000
00180000
00190000
00200000
00210000
00220000
00230001
00240001
00250001
00260001
00270001
00280001
00290000
00300000
00310000
00320000
00330000
00340001
00350001

```

```

+ /'00','@','$','#','Z' '/'
C*-----< 6 LINES OMIT (93/05/17)
C+ INTEGER MICKOD(34)
C+ 1 /'1','2','3','4','5','6','7','8',
C+ 2 '9','A','B','C','D','E','F','G',
C+ 3 'H','I','J','K','L','M','N','O',
C+ 4 'P','Q','R','S','T','U','V','W',
C+ 5 'X','Y' '/'
C
C IF (KACC.LT.1 .OR. KACC.GT.6) GO TO 7000
C
C GO TO (1100,1200,1300,1400,1500,1600) ,KACC
C
C MEMBER NAME FOR ID DATA
C
1100 CONTINUE
ID = IDCOD
GO TO 2000
C
C MEMBER NAME FOR MACRO DATA
C
1200 CONTINUE
ID = MACRO
GO TO 2000
C
C MEMBER NAME FOR MICRO DATA
C
1300 CONTINUE
IF (NMIC .LE. 0 ) GO TO 7100
C*-----< 1 LINE OMIT (93/05/17)
C* ID = MICKOD(1)
DO 1310 I=1,NMIC
IF (MICCOD.NE.NMICRO(I)) GO TO 1310
C*-----< 1 LINE OMIT & 2 LINES INSERT (93/05/17)
C+ ID = MICKOD(I)
WRITE(NUM,'(2X,I2)') I
IF ( NUM(3:3).EQ.' ' ) NUM(3:3) = '0'
READ(NUM,'(2X,A2)') ID
C*
GO TO 1320
1310 CONTINUE
GO TO 7100
1320 CONTINUE
GO TO 2000
C
C MEMBER NAME FOR BUCKLING DATA
C
1400 CONTINUE
ID = KBUCK
GO TO 2000
C
C MEMBER NAME FOR FISSION SPECTRUM DATA
C
1500 CONTINUE
ID = KHI
GO TO 2000
C
C MEMBER NAME FOR ZONE AVERAGED FLUX DATA
C
1600 CONTINUE
ID = LUXF
C
C
2000 CONTINUE
C
C
ID1 = 10
DO 100 I=1,10
IF ( KODE(I,1).NE.NMPROG ) GO TO 100
ID1 = I
GO TO 200

```

```

00360001
00370001
00380001
00390001
00400001
00410001
00420001
00430001
00440000
00450000
00460000
00470000
00480000
00490000
00500000
00510000
00520000
00530000
00540000
00550000
00560000
00570000
00580000
00590000
00600000
00610001
00620001
00630000
00640000
00650001
00660001
00670001
00680001
00690001
00700001
00710000
00720000
00730000
00740000
00750000
00760000
00770000
00780000
00790000
00800000
00810000
00820000
00830000
00840000
00850000
00860000
00870000
00880000
00890000
00900000
00910000
00920000
00930000
00940000

```



```
//FT26F001 DD DSN=&KEFFC,UNIT=SYSDA,SPACE=(TRK,(10,10)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT27F001 DD DSN=&RPEAK,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT28F001 DD DSN=&POFLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT31F001 DD DSN=&AJFLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT32F001 DD DSN=&FTFLY,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT39F001 DD DSN=&P13PO,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT40F001 DD DSN=&P6FLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT42F001 DD DSN=&DLLEAK,UNIT=SYSDA,SPACE=(TRK,(10,10)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT44F001 DD DSN=&APLUX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT45F001 DD DSN=&APQWR,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT46F001 DD DSN=&P6POW,UNIT=SYSDA,SPACE=(TRK,(100,25)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT50F001 DD DSN=&AVTEM,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT54F001 DD DSN=&TDEPN,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT55F001 DD DSN=&FLUX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT56F001 DD DSN=&FTAU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT57F001 DD DSN=&SOURC,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT58F001 DD DSN=&PTSAE,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT59F001 DD DSN=&DLEK1,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT60F001 DD DSN=&FTFLX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT61F001 DD DSN=&P6HTX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT16F001 DD DSN=&HIST1,UNIT=SYSDA,SPACE=(TRK,(5,5)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT17F001 DD DSN=&HIST2,UNIT=SYSDA,SPACE=(TRK,(5,5)),DISP=(,PASS),
//          DCB=*.FT12F001
//PLOTLOG DD DUMMY
//GDFILE DD DUMMY
//*PLOTLOG DD SYSOUT=&OUT,DCB=*.FT06F001
//*GDFILE DD SYSOUT=&OUT,DCB=*.FT06F001
//*PLOTPRM DD *
//* SCALE=0.7
//-----
//FT01F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=*.FT12F001
//FT02F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(&N4)),DCB=*.FT01F001
//FT15F001 DD UNIT=SYSDA,SPACE=(TRK,(&N7)),DCB=*.FT01F001
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(&N8)),DCB=*.FT15F001
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT19F001 DD UNIT=SYSDA,SPACE=(TRK,(&N9)),DCB=*.FT01F001
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(&N10)),DCB=*.FT01F001
```

```
//FT34F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT35F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=*.FT01F001
//FT36F001 DD UNIT=SYSDA,SPACE=(TRK,(&N11)),DCB=*.FT01F001
//FT37F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=*.FT01F001
//FT38F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT43F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT51F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT52F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=*.FT01F001
//FT62F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//FT63F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=*.FT01F001
//*--- INTERFACE FILE
//FT74F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//FT75F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//FT76F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//FT77F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//FT79F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//FT82F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=*.FT01F001
//          PEND
//*
//          EXEC MOSESV1,
//          N1=100,N2=6526,N3=10,N4=1,N5=2877,N6=10,
//          N7=103,N8=103,N9=933,N10=933,N11=1015,N12=1,N13=1,N14=1,
//          N15=10,N16=6526,NX=60,NS=1015
//PLOTLOG DD DUMMY
//GDFILE DD SYSOUT=R
//PLOTPRM DD *
//          SCALE=0.7
//*
//FT03F001 DD DISP=SHR,DSN=POC0B17.OBHZHBC7.MIC7GJ3,LABEL=(,,IN)
//FT05F001 DD DISP=SHR,LABEL=(,,IN),
//          DSN=POC0B17.OPMOSD.DATA(GDI1DY)
//
```

JCL file [20] .OPJCL.CNTL(GDI1KSLA)

SLAROM, 'shutdown margin'

```
//POC0B17X JOB (),'PTSLAROM',MSGCLASS=X,NOTIFY=POC0B17,MSGLEVEL=(2,0), 00010042
// CLASS=B,TIME=0005 00020001
//***** 00030000
//*** SLAROM *** FOR PERTURBATION, INC.DOPP.-DENSITY 00040029
//***** 00050000
//DATAP EXEC PGM=DATAP 00060032
//STEPLIB DD DSN=POC0H##.DATAP.LOAD,DISP=SHR,LABEL=(,,IN) 00070032
//FT06F001 DD SYSOUT=* 00080014
//* ++++++ INPUT SLAROM N.D.(FUEL*3) 00110021
//FT05F001 DD DSN=POC0B17.OPPEND.DATA(GDI1RZEC),DISP=SHR, 00120045
//* ----- 00121002
//          LABEL=(,,IN) 00130000
//FT09F001 DD UNIT=SYSDA,DSN=&&INPUT,DISP=(,PASS),SPACE=(CYL,(1,1)) 00131014
//***** 00140000
//GO EXEC PGM=SLAROM 00160033
//STEPLIB DD DSN=POC0EAL.SLAROM.NEDAC.LOAD,DISP=SHR 00180033
//*----- ( INPUT PDS )----- 00190000
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//PDSIN DD DUMMY 00200000
//* ***** OUTPUT 70 GROUP PDS 00210021
//PDSOUT DD DSN=POC0B17.@GDI1P.PDS70G,DISP=SHR 00220046
//* 00221021
//* DISP=(NEW,CATLG),UNIT=DASD, 00230046
//* DCB=(RECFM=FB,BLKSIZE=4000,LRECL=4,DSORG=PO), 00240046
//* SPACE=(CYL,(10,20,50),RLSE) 00250046
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)), 00270000
// DCB=(RECFM=VSB,LRECL=32756,BLKSIZE=32760) 00280015
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00290000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00300000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00310000
//FT05F001 DD DDNAME=SYSIN 00330014
//FT06F001 DD SYSOUT=*,DCB=(RECFM=FB,LRECL=137,BLKSIZE=1370) 00350014
//FT06F002 DD SYSOUT=*,DCB=(RECFM=FB,LRECL=137,BLKSIZE=1370) 00371014
//FT07F001 DD DUMMY 00380000
//* ***** INPUT REF. 70 GROUP MICRO00390002
//FT08F001 DD DSN=POCOH#.JFS3J3.Y9406,DISP=SHR,LABEL=(,IN) 00420010
//* 00430002
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(10,10)),DCB=*.FT01F001 00440014
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00450000
//FT11F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00460000
//FT12F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00470000
//FT13F001 DD DUMMY 00471014
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(10,5)),DCB=*.FT01F001 00480000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00490000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00500000
//FT22F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00510000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00520000
//FT26F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00530000
//FT30F001 DD DUMMY 00540000
//FT40F001 DD SYSOUT=* 00550000
//FT41F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00560000
//FT42F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00570000
//FT43F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00580000
//FT55F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00590000
//FT56F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00600000
//FT59F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00610000
//FT60F001 DD SYSOUT=* 00620014
//FT70F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00630000
//FT72F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00640000
//FT73F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00650000
//FT71F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00660000
//FT80F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00670000
//FT90F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00680000
//FT96F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00690000
//FT99F001 DD UNIT=SYSDA,SPACE=(CYL,(5,5)),DCB=*.FT01F001 00700000
//SYSIN DD DSN=*&INPUT,DISP=(OLD,DELETE) 00730015
// 02410000

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JCL file [21] .OPJCL.CNTL(GDI1X70N)

CITATION, normal, 70 groups, 'shutdown margin'

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//POC0B17X JOB (),'NEMC7GOR',MSGCLASS=X,NOTIFY=POC0B17,MSGLEVEL=(2,0), 00010042
// CLASS=B,TIME=0005 00020001
//***** 000770000
//* JOINT ---> CITATION **//00790000
//***** 00810000
//JOINT1 EXEC PGM=JOINTX 00950032
//STEPLIB DD DSN=POC0EAL.JOINT.LOAD,DISP=SHR 00960032
//SYSPRINT DD SYSOUT=* 00970000
//* ***** INPUT 70GP. MIC&MAC. 00971015
//USERPDS DD DSN=POC0B17.@GDI1P.PDS70G,DISP=SHR 00980044
//* 00981015
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,10)), 00991031
// DISP=(,PASS),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 01000031
//* ***** INPUT CIT.DATA 70GP/D&A 01010019
//FT05F001 DD DSN=POC0B17.OPCITD.DATA(PFC3N70X),DISP=SHR, 01011038
//* 01012002
// LABEL=(,IN) 01020000
//FT06F001 DD SYSOUT=* 01030031
//*FT06F001 DD DSN=*&LIST4,DISP=(NEW,PASS), 01031031
//* SPACE=(CYL,(10,3)), 01040031
//* DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 01050031
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01060000
// SPACE=(TRK,(10,10)), 01070000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01080000
//FT09F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01090000
// SPACE=(TRK,(10,10)) 01100000
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01110000
// SPACE=(TRK,(10,10)), 01120000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01130000
//FT20F001 DD DUMMY 01140000
//FT50F001 DD SYSOUT=* 01150000
//* 01160000
//*L0UT4 EXEC COMPACT,PARM='TYPE2,CC=NO' 01170031
//*UTYIN DD DSN=*&LIST4,DISP=(OLD,DELETE) 01180031
//*----- CITATION ----- 01190000
//CITATION EXEC PGM=CITATION,COND=(4,LT) 01200000
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR 01210000
//SYSPRINT DD SYSOUT=* 01220000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01230000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 01240000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01250000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01260000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01270000
//FT05F001 DD DSN=*.JOINT1.FT08F001,DISP=(OLD,DELETE) 01280000
//FT06F001 DD SYSOUT=* 01290000
//*---( OUTPUT ZONE AVERAGED FLUX & AXIAL BUCKLING )-----** 01300000
//FT07F001 DD DSN=*&CIT,DISP=(NEW,PASS),UNIT=WORK, 01310000
// SPACE=(TRK,(2,1),RLSE), 01320000
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200) 01330000
//FT08F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01340000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 01350000
//*---( OUTPUT FLUX MAP )-----** 01360000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01370000

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//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01380000 //PDSIN DD DSN=POCOB17. @GDI1P.PDS70G,DISP=SHR 01930044
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01390000 //** 01931015
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01400000 //** ***** OUTPUT 18GP. PDS MIC&MAC01940018
//**---( RESTART FILE )-----** 01410000 //PDSOUT DD DSN=POCOB17. @GDI1P18. PDSMICN, 01960044
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01420000 //** 01961002
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01430000 // UNIT=DASD,DCB=(RECFM=FB,LRECL=4,BLKSIZE=4000),DISP=(NEW,CATLG), 01970000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(2,5)),DCB=*.FT01F001 01440000 // SPACE=(TRK,(7,1,30)) 01980021
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01450000 //** ***** OUTPUT 70 GROUP FLUX01990002
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01460000 //USERPDS DD DSN=POCOB17. @GDI1P70. FLUXN,DISP=SHR 02000044
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01470000 //** 02001002
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01480000 //FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 02010000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01490000 //** ***** INPUT CONDENSATION CONTROL02020002
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01500000 //FT05F001 DD DSN=POCOB17. OPEDID. DATA (PFC3C18X),DISP=SHR, 02021038
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01510000 //** 02022002
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01520000 // LABEL=(, , , IN) 02030000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01530000 //FT06F001 DD SYSOUT=* 02040000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01540000 //FT90S001 DD SYSOUT=* 02050000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01550000 //**---( JOINT )----- 02060025
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01560000 //MICRO EXEC PGM=JOINTX 02070032
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01570000 //STEPLIB DD DSN=POCOEAL. JOINT. LOAD,DISP=SHR,LABEL=(, , , IN) 02080032
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01580000 //SYSPRINT DD SYSOUT=* 02090025
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01590000 // DCB=(RECFM=FBA,LRECL=137,BLKSIZE=19043) 02100025
//**---( MACRO CROSS SECTIONS )----- 01600000 //** ***** OUTPUT 7 GP. MICRO (PDS)02110025
//FT31F001 DD DSN=*.JOINT1. FT09F001,DISP=(OLD,DELETE) 01610000 //USERPDS DD DSN=POCOB17. @GDI1P18. PDSMICN,DISP=SHR 02120044
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01620000 //** 02130025
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01630000 //FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)), 02140025
//** ***** OUTPUT 70 GP FLUX DIR&ADJ. 01640019 // DISP=(, PASS),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 02150025
//FT34F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01650026 //** ***** INPUT MICRO EDIT CONTROL02160025
//FT34F001 DD DSN=POCOB17. @LCM4P. FT34G70N,DISP=(NEW,CATLG), 01650126 //FT05F001 DD DSN=POCOB17. OPEDID. DATA (PFC3MX),DISP=SHR 02170038
//** 01651019 //** 02180025
//** UNIT=DASD,SPACE=(TRK,(5,5),RLSE),DCB=*.FT01F001 01652030 //FT06F001 DD SYSOUT=* 02190031
//FT51F001 DD SYSOUT=* 01660031 //FT06F001 DD DSN=&&LIST6,DISP=(NEW,PASS), 02191031
//FT51F001 DD DSN=&&LIST5,DISP=(NEW,PASS), 01661031 //** SPACE=(CYL,(10,3)), 02200031
//** 01670031 //** DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 02210031
//** 01680031 //FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 02220025
//** 01690000 // SPACE=(TRK,(10,10)), 02230025
//** 01700031 // DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 02240025
//** 01710031 //** ***** OUTPUT 7 GP. MICRO (SLAROM)02250025
//**---( JOINT.MIC )----- 01720000 //FT09F001 DD DSN=POCOB17. @GDI1P18. MICRON,DISP=(NEW,CATLG),UNIT=SYSDA, 02260044
//JOINT2 EXEC PGM=JOINTX 01750032 //** 02270025
//STEPLIB DD DSN=POCOEAL. JOINT. LOAD,DISP=SHR,LABEL=(, , , IN) 01760033 // SPACE=(TRK,(10,5),RLSE), 02280025
//SYSPRINT DD SYSOUT=* 01770000 // DCB=(RECFM=VBS,LRECL=X,BLKSIZE=26793) 02290025
//** ***** OUTPUT 70 GROUP FLUX01780002 //FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 02300025
//USERPDS DD DSN=POCOB17. @GDI1P70. FLUXN,DISP=(NEW,CATLG),UNIT=DASD, 01781044 // SPACE=(TRK,(10,10)), 02310025
//** 01782002 // DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 02320025
//** 01790000 //FT20F001 DD DUMMY 02330025
//** 01800000 //FT50F001 DD SYSOUT=* 02340025
//FT01F001 DD DSN=&&CIT,DISP=(OLD,DELETE) 01810000 //** 02350025
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 01820000 //** 02360025
//** ***** INPUT FLUX EDIT CONTROL01830002 //** 02370031
//FT05F001 DD DSN=POCOB17. OPEDID. DATA (PFC3FX),DISP=SHR 01831038 //LOUT6 EXEC COMPACT,PARM='TYPE2,CC=NO' 02380031
//** 01832002 //UTYIN DD DSN=&&LIST6,DISP=(OLD,DELETE) 02410000
//** 01840000 //** 02410000
//FT90S001 DD SYSOUT=* 01850000
//**---( JOINT.MIC )----- 01870000
//JOINT3 EXEC PGM=JOINTX 01900032
//STEPLIB DD DSN=POCOEAL. JOINT. LOAD,DISP=SHR,LABEL=(, , , IN) 01910032
//SYSPRINT DD SYSOUT=* 01920000
//** ***** INPUT 70GP. PDS (MIC&MAC) 01921015

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JCL file [22] .OPJCL.CNTL(GDI1X70D)

CITATION, Doppler, 70 groups, 'shutdown margin'

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//POCOB17X JOB (),'NEMC7GOR',MSGCLASS=X,NOTIFY=POCOB17,MSGLEVEL=(2,0), 00010039
// CLASS=B,TIME=0005 00020001
//*****//00770000
//* JOINT --> CITATION *//00790000
//*****//00810000
//JOINT1 EXEC PGM=JOINTX 00950033
//STEPLIB DD DSN=POCOEAL.JOINT.LOAD,DISP=SHR 00960033
//SYSPRINT DD SYSOUT=* 00970000
//* ***** INPUT 70GP. MIC&MAC. 00971015
//USERPDS DD DSN=POCOB17. @GDI1P. PDS70G,DISP=SHR 00980041
//* ***** 00981015
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,10)), 00990000
// DISP=(,PASS),DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200,BUFNO=1) 01000000
//* ***** INPUT CIT. DATA 70GP/D&A 01010019
//FT05F001 DD DSN=POCOB17. OPCITD. DATA (PFC3N70D),DISP=SHR, 01011036
//* ***** 01012002
// LABEL=(, , IN) 01020000
//FT06F001 DD SYSOUT=* 01030033
//*FT06F001 DD DSN=&&LIST4,DISP=(NEW,PASS), 01031033
//* SPACE=(CYL,(10,3)), 01040033
//* DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 01050033
//FT08F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01060000
// SPACE=(TRK,(10,10)), 01070000
//* DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01080000
//FT09F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01090000
// SPACE=(TRK,(10,10)) 01100000
//FT10F001 DD DISP=(NEW,PASS,DELETE),UNIT=SYSDA, 01110000
// SPACE=(TRK,(10,10)), 01120000
//* DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120) 01130000
//FT20F001 DD DUMMY 01140000
//FT50F001 DD SYSOUT=* 01150000
//* 01160000
//*LOUT4 EXEC COMPACT,PARM='TYPE2,CC=NO' 01170033
//*UTYIN DD DSN=&&LIST4,DISP=(OLD,DELETE) 01180033
//*----- ( CITATION )----- 01190000
//CITATION EXEC PGM=CITATION,COND=(4,LT) 01200000
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR 01210000
//SYSPRINT DD SYSOUT=* 01220000
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01230000
// DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 01240000
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01250000
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01260000
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01270000
//FT05F001 DD DSN=*.JOINT1.FT08F001,DISP=(OLD,DELETE) 01280000
//FT06F001 DD SYSOUT=* 01290000
//*----- ( OUTPUT ZONE AVERAGED FLUX & AXIAL BUCKLING )----- 01300000
//FT07F001 DD DSN=&&CIT,DISP=(NEW,PASS),UNIT=WORK, 01310000
// SPACE=(TRK,(2,1),RLSE), 01320000
//* DCB=(RECFM=FB,LRECL=80,BLKSIZE=3200) 01330000
//FT08F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)), 01340000
//* DCB=(RECFM=VSB,LRECL=4092,BLKSIZE=4096) 01350000
//*----- ( OUTPUT FLUX MAP )----- 01360000
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01370000

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//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01380000
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01390000
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01400000
//*----- ( RESTART FILE )----- 01410000
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01420000
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01430000
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(2,5)),DCB=*.FT01F001 01440000
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01450000
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01460000
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(50,5)),DCB=*.FT01F001 01470000
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01480000
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01490000
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01500000
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01510000
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01520000
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01530000
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01540000
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01550000
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01560000
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01570000
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01580000
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=*.FT01F001 01590000
//*----- ( MACRO CROSS SECTIONS )----- 01600000
//FT31F001 DD DSN=*.JOINT1.FT09F001,DISP=(OLD,DELETE) 01610000
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01620000
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001 01630000
//* ***** OUTPUT 70 GP FLUX DIR&ADJ. 01640019
//FT34F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001 01650026
//*FT34F001 DD DSN=POCOB17. @LCM4P. FT34G70N,DISP=(NEW,CATLG), 01650126
//* ***** 01651019
//* UNIT=DASD,SPACE=(TRK,(5,5),RLSE),DCB=*.FT01F001 01652030
//FT51F001 DD SYSOUT=* 01660033
//*FT51F001 DD DSN=&&LIST5,DISP=(NEW,PASS), 01661033
//* SPACE=(CYL,(10,3)), 01670033
//* DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK 01680033
//* 01690000
//*LOUT5 EXEC COMPACT,PARM='TYPE2,CC=YES' 01700033
//*UTYIN DD DSN=&&LIST5,DISP=(OLD,DELETE) 01710033
//*----- ( JOINT.MIC )----- 01720000
//JOINT2 EXEC PGM=JOINTX 01750033
//STEPLIB DD DSN=POCOEAL.JOINT.LOAD,DISP=SHR,LABEL=(, , IN) 01760033
//SYSPRINT DD SYSOUT=* 01770000
//* ***** OUTPUT 70 GROUP FLUX 01780002
//USERPDS DD DSN=POCOB17. @GDI1P70. FLUXD,DISP=(NEW,CATLG),UNIT=DASD, 01781041
//* ***** 01782002
//* DCB=(RECFM=FB,LRECL=4,BLKSIZE=4000), 01790000
//* SPACE=(TRK,(5,1,3),RLSE) 01800000
//FT01F001 DD DSN=&&CIT,DISP=(OLD,DELETE) 01810000
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 01820000
//* ***** INPUT FLUX EDIT CONTROL 01830002
//FT05F001 DD DSN=POCOB17. OPEDID. DATA (PFC3FD),DISP=SHR 01831036
//* ***** 01832002
//FT06F001 DD SYSOUT=* 01840000
//FT90S001 DD SYSOUT=* 01850000
//*----- ( JOINT.MIC )----- 01870000
//JOINT3 EXEC PGM=JOINTX 01900033
//STEPLIB DD DSN=POCOEAL.JOINT.LOAD,DISP=SHR,LABEL=(, , IN) 01910033
//SYSPRINT DD SYSOUT=* 01920000
//* ***** INPUT 70GP. PDS (MIC&MAC) 01921015

```

```

//PDSIN DD DSN=POC0B17.@GDI1P.PDS70G,DISP=SHR 01930041
//* 01931015
//* ***** OUTPUT 18GP. PDS MIC&MAC01940018
//PDSOUT DD DSN=POC0B17.@GDI1P18.PDSMICD, 01960041
//* 01961002
// UNIT=DASD,DCB=(RECFM=FB,LRECL=4,BLKSIZE=4000),DISP=(NEW,CATLG), 01970000
// SPACE=(TRK,(7,1,30)) 01980021
//* ***** OUTPUT 70 GROUP FLUX01990002
//USERPDS DD DSN=POC0B17.@GDI1P70.FLUXD,DISP=SHR 02000041
//* 02001002
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(10,2)) 02010000
//* ***** INPUT CONDENSATION CONTROL02020002
//FT05F001 DD DSN=POC0B17.OPEDID.DATA(PFC3C18D),DISP=SHR, 02021037
//* 02022002
// LABEL=(, , , IN) 02030000
//FT06F001 DD SYSOUT=* 02040000
//FT90S001 DD SYSOUT=* 02050000
//*****-( JOINT )-----02060025
// 02410000

```

JCL file [23] .OPJCL.CNTL(GDI1PERK)

PERKY package, 'shutdown margin'

```

//POC0B17B JOB (,'FBRJ',MSGCLASS=X,NOTIFY=POC0B17,
// MSGLEVEL=(1,1),CLASS=2,TIME=0010
//* SHORTENED , MODIFIED TO SNH FILENAMES
//*****
//** CASE PIJ1 - FINAL OPTIMIZED 165% LOQ 45.05% 4*10.5M 21/OP ZRH
//*****
//***** PRE-PERKY
//PREPK EXEC PGM=PREPK99
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR
//***** N.D. FILE *****
//FT01F001 DD DSN=POC0B17.@GDI1.DENS,DISP=SHR
//*****
//* !!
//***** CITATION INPUT, FOR N.D. *****
//FT02F001 DD DSN=POC0B17.@GDI1.CITDAT,DISP=SHR,LABEL=(, , , IN)
//*****
//FT03F001 DD DSN=POC0B17.OPEDID.DATA(NAMEF#XX), INPUT
// DISP=SHR,LABEL=(, , , IN) INPUT
//FT04F001 DD DSN=POC0B17.OPEDID.DATA(IDNUM#XX), INPUT
// DISP=SHR,LABEL=(, , , IN) INPUT
//FT14F001 DD DSN=POC0B17.OPEDID.DATA(CHIMG), INPUT
// DISP=SHR,LABEL=(, , , IN) INPUT
//*
//** CYCLE STEP SELECT-STEP
//FT05F001 DD * INPUT
// 6 2 2
//*
//*

```

```

//** 18G PDS BASE FILE (FT16F001)
//** 18G PDS BASE+500C FILE (FT17F001)
//**
//***** NORMAL PDS FILE *****
//FT16F001 DD DSN=POC0B17.@GDI1P18.PDSMICN, INPUT
//*****
// DISP=SHR,LABEL=(, , , IN) INPUT
//***** DOPPLER PDS FILE *****
//FT17F001 DD DSN=POC0B17.@GDI1P18.PDSMICD, INPUT
//*****
// DISP=SHR,LABEL=(, , , IN) INPUT
//**
//**
//FT07F001 DD DSN=POC0B17.@CITE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT21F001 DD DSN=POC0B17.@DOPE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT22F001 DD DSN=POC0B17.@FULE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT23F001 DD DSN=POC0B17.@STRE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT24F001 DD DSN=POC0B17.@COLE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT25F001 DD DSN=POC0B17.@VIDE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT26F001 DD DSN=POC0B17.@BETE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT27F001 DD DSN=POC0B17.@LIFE,SPACE=(TRK,(1,1)),UNIT=DASD,
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120),DISP=(NEW,CATLG)
//FT31F001 DD DSN=POC0B17.@PMICE,SPACE=(TRK,(20,20)),UNIT=DASD,
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096),DISP=(NEW,CATLG)
//FT90F001 DD DUMMY
//FT06F001 DD SYSOUT=*
//**FT06F001 DD DSN=&&DALIS0,DISP=(NEW,DELETE),
//* SPACE=(TRK,(100,50)),
//** DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//**
//***** CITATION ADJOINT
//ADJOINT EXEC PGM=CITATION,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.CITFBR.LOAD,DISP=SHR
//FT01F001 DD UNIT=SYSDA,SPACE=(CYL,(20,2)),
// DCB=(RECFM=VSE,LRECL=4092,BLKSIZE=4096)
//FT02F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=* FT01F001
//FT03F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=* FT01F001
//FT04F001 DD UNIT=SYSDA,SPACE=(CYL,(1,2)),DCB=* FT01F001
//*****
//FT05F001 DD DSN=POC0B17.@CITE,DISP=SHR
//*****
//FT06F001 DD DSN=&&DALIST,DISP=(NEW,PASS),
// SPACE=(TRK,(30,20)),
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//FT07F001 DD UNIT=SYSDA,SPACE=(TRK,(50,50)),DCB=* FT01F001
//*****
//** ! CITATION 18G MICRO X-SECTION ! **
//FT08F001 DD DSN=POC0B17.@GDI1P18.MICRON,DISP=SHR,LABEL=(, , , IN) INPUT
//*****

```

```

/**
//FT09F001 DD UNIT=SYSDA,SPACE=(CYL,(15,8)),DCB=*.FT01F001
//FT10F001 DD UNIT=SYSDA,SPACE=(CYL,(4,2)),DCB=*.FT01F001
//FT11F001 DD UNIT=SYSDA,SPACE=(TRK,(100,5)),DCB=*.FT01F001
//FT12F001 DD UNIT=SYSDA,SPACE=(TRK,(100,5)),DCB=*.FT01F001
//FT13F001 DD UNIT=SYSDA,SPACE=(TRK,(100,5)),DCB=*.FT01F001
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(100,5)),DCB=*.FT01F001
//FT15F001 DD UNIT=SYSDA,SPACE=(CYL,(8,5)),DCB=*.FT01F001
//FT16F001 DD UNIT=SYSDA,SPACE=(TRK,(80,5)),DCB=*.FT01F001
//FT17F001 DD UNIT=SYSDA,SPACE=(TRK,(80,5)),DCB=*.FT01F001
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(80,5)),DCB=*.FT01F001
//FT19F001 DD UNIT=SYSDA,SPACE=(CYL,(20,8)),DCB=*.FT01F001
//FT20F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT21F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT22F001 DD UNIT=SYSDA,SPACE=(TRK,(20,10)),DCB=*.FT01F001
//FT23F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT24F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT25F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT26F001 DD UNIT=SYSDA,SPACE=(TRK,(20,10)),DCB=*.FT01F001
//FT27F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT28F001 DD UNIT=SYSDA,SPACE=(TRK,(20,10)),DCB=*.FT01F001
//FT29F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
//FT30F001 DD UNIT=SYSDA,SPACE=(CYL,(2,2)),DCB=*.FT01F001
/*-----
//FT31F001 DD DSN=POCOB17.@PMACE,UNIT=DASD,
// DISP=(NEW,CATLG),SPACE=(TRK,(6,2),RLSE),
// DCB=(RECFM=FB,LRECL=80,BLKSIZE=3120)
//FT34F001 DD DSN=POCOB17.@FT34E,UNIT=DASD,
// DISP=(NEW,CATLG),SPACE=(TRK,(60,20),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
/*-----
//FT32F001 DD UNIT=SYSDA,SPACE=(TRK,(10,10)),DCB=*.FT01F001
//FT33F001 DD UNIT=SYSDA,SPACE=(CYL,(8,8)),DCB=*.FT01F001
//FT51F001 DD SYSOUT=*
//FT51F001 DD DSN=*&LTCIT,DISP=(NEW,PASS),
// SPACE=(CYL,(15,3)),
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//FT52F001 DD UNIT=SYSDA,SPACE=(CYL,(16,8)),DCB=*.FT01F001
//FT61F001 DD DUMMY
//FT80F001 DD DUMMY
//FT81F001 DD DUMMY
/**
***** PERKY DOPPLER
***** PERKY FUEL DENS
//PERFUL EXEC PGM=PERKY,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR,LABEL=(, , IN)
//SYSPRINT DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/*-----
//FT01F001 DD DSN=POCOB17.@FT34E,DISP=SHR
/*-----
//FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DISP=(NEW,DELETE),
// DCB=(RECFM=VBS,LRECL=23000,BLKSIZE=23004)
//FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*-----
//FT05F001 DD DSN=POCOB17.@FULE,DISP=(OLD,PASS)

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```

/*-----
//FT06F001 DD DSN=*&FULOUT,DISP=(NEW,DELETE),
// SPACE=(TRK,(100,10)),
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//FT08F001 DD DUMMY
/*MACRO1
//FT09F001 DD DUMMY
/*MACRO2
//FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*-----
//FT11F001 DD DSN=POCOB17.@PMICE,DISP=SHR
//FT20F001 DD DSN=POCOB17.@PMACE,DISP=SHR
/*-----
//FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT21F001 DD DSN=POCOB17.DELAY.DATA,LABEL=(, , IN),DISP=SHR
//FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT25F001 DD DSN=*&PERK25,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT27F001 DD DSN=*&PERK27,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/**
//POSFUL EXEC PGM=POSTPK,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR
//SYSPRINT DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/** INPUT
//FT05F001 DD DSN=POCOB17.PQINT.DATA(POFUL),
// DISP=SHR,LABEL=(, , IN)
INPUT
INPUT
/**
//FT06F001 DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425)
/** INPUT
//FT10F001 DD DSN=POCOB17.OPEDID.DATA(ZONE#XX),
// DISP=SHR,LABEL=(, , IN)
INPUT
INPUT
/**
//FT25F001 DD DSN=*&PERK25,DISP=(OLD,DELETE)
//FT27F001 DD DSN=*&PERK27,DISP=(OLD,DELETE)
/*-----
//FT34F001 DD DSN=POCOB17.@FT34E,DISP=SHR
//FT55F001 DD DSN=POCOB17.@CITE,DISP=SHR
/*-----
/**
***** PERKY STRUCTURE
//PERSTR EXEC PGM=PERKY,COND=(4,LT)

```

```

//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR,LABEL=(, , , IN)
//SYSPRINT DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/*-----
//FT01F001 DD DSN=POC0B17.@FT34E,DISP=SHR
//FT05F001 DD DSN=POC0B17.@STRE,DISP=SHR
/*-----
//FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DISP=(NEW,DELETE),
// DCB=(RECFM=VBS,LRECL=23000,BLKSIZE=23004)
//FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT06F001 DD DSN=&&STROUT,DISP=(NEW,DELETE),
// SPACE=(TRK,(100,20)),
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//FT08F001 DD DUMMY
/*MACRO1
//FT09F001 DD DUMMY
/*MACRO2
//FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*-----
//FT11F001 DD DSN=POC0B17.@PMICE,DISP=SHR
//FT20F001 DD DSN=POC0B17.@PMACE,DISP=SHR
/*-----
//FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT21F001 DD DSN=POC0B17.DELAY.DATA,LABEL=(, , , IN),DISP=SHR
//FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT25F001 DD DSN=&&PERK25,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT27F001 DD DSN=&&PERK27,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*
/*
//POSTR EXEC PGM=POSTPK,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR
//SYSPRINT DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/*
/* INPUT
//FT05F001 DD DSN=POC0B17.PQINT.DATA(POSTR), INPUT
// DISP=SHR,LABEL=(, , , IN) INPUT
/*
//FT06F001 DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425)
/*
/* INPUT
//FT10F001 DD DSN=POC0B17.OPEDID.DATA(ZONE#XX), INPUT

```

```

// DISP=SHR,LABEL=(, , , IN)
/*
//FT25F001 DD DSN=&&PERK25,DISP=(OLD,DELETE)
//FT27F001 DD DSN=&&PERK27,DISP=(OLD,DELETE)
/*-----
//FT34F001 DD DSN=POC0B17.@FT34E,DISP=SHR
//FT55F001 DD DSN=POC0B17.@CITE,DISP=SHR
/*-----
/*
***** PERKY COOLANT
//PERCOL EXEC PGM=PERKY,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR,LABEL=(, , , IN)
//SYSPRINT DD SYSOUT=*,
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/*-----
//FT01F001 DD DSN=POC0B17.@FT34E,DISP=SHR
//FT05F001 DD DSN=POC0B17.@COLE,DISP=SHR
/*-----
//FT02F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DISP=(NEW,DELETE),
// DCB=(RECFM=VBS,LRECL=23000,BLKSIZE=23004)
//FT03F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT04F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT06F001 DD DSN=&&COLOUT,DISP=(NEW,DELETE),
// SPACE=(TRK,(100,50)),
// DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425),UNIT=WORK
//FT08F001 DD DUMMY
/*MACRO1
//FT09F001 DD DUMMY
/*MACRO2
//FT10F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*-----
//FT11F001 DD DSN=POC0B17.@PMICE,DISP=(OLD,PASS)
//FT20F001 DD DSN=POC0B17.@PMACE,DISP=(OLD,PASS)
/*-----
//FT13F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT14F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT15F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT16F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT18F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT19F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT21F001 DD DSN=POC0B17.DELAY.DATA,LABEL=(, , , IN),DISP=SHR
//FT22F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT25F001 DD DSN=&&PERK25,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT27F001 DD DSN=&&PERK27,UNIT=WORK,DISP=(NEW,PASS),
// SPACE=(TRK,(10,2),RLSE),
// DCB=(RECFM=VBS,LRECL=4092,BLKSIZE=4096)
//FT31F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT32F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT41F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
//FT42F001 DD UNIT=WORK,SPACE=(TRK,(100,50)),DCB=*.FT02F001
/*
/*
//POSCOL EXEC PGM=POSTPK,COND=(4,LT)
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DSN=POC3A12.PERKYL2.CS.LOAD,DISP=SHR
//SYSPRINT DD SYSOUT=*,

```

INPUT

```

//          DCB=(RECFM=FBA,LRECL=137,BLKSIZE=8220)
/**
/** INPUT
//FT05F001 DD DSN=POCOB17.PQINT.DATA(POCOL),
//          DISP=SHR,LABEL=(,,,IN)
/**
//FT06F001 DD SYSOUT=*,
//          DCB=(RECFM=FBA,LRECL=137,BLKSIZE=3425)
/**
/** INPUT
//FT10F001 DD DSN=POCOB17.OPEDID.DATA(ZONE#XK),
//          DISP=SHR,LABEL=(,,,IN)
/**
//FT25F001 DD DSN=&&PERK25,DISP=(OLD,DELETE)
//FT27F001 DD DSN=&&PERK27,DISP=(OLD,DELETE)
/**-----
//FT34F001 DD DSN=POCOB17.@FT34E,DISP=SHR
//FT55F001 DD DSN=POCOB17.@CITE,DISP=SHR
/**-----
/**
//***** PERKY VOID
//***** PERKY BETA-BFF
//***** PERKY LIFE-TIME
/**
//

```

INPUT INPUT INPUT INPUT

JCL file [24] .OPJCL.CNTL(GDI1MBRN)
MOSES burn-up, '3D burn-up'

```

//POCOB17Z JOB (),MOSESV1,NOTIFY=POCOB17,
//          MSGCLASS=X,MSGLEVEL=(1,1),CLASS=4,TIME=0120
/**
//MOSESV1 PROC MODULE=MOSES96,OUT='*',STEPDS='#MOSES96.LOAD',
//          NK=200,NS=5000,N1=100,
//          N4=1,N7=103,N8=103,N9=933,N10=2414,N11=1015
//GO EXEC PGM=&MODULE,PARM='FLIB(RUN77)'
//VPSYS DD SUBSYS=(VPCS)
//STEPLIB DD DISP=SHR,DSN=POC3A12.&STEPDS
//PRINT DD SYSOUT=&OUT,DCB=(RECFM=FBA,LRECL=133,BLKSIZE=1330)
//FT06F001 DD SYSOUT=&OUT,DCB=(RECFM=FBA,LRECL=133,BLKSIZE=1330)
/**FT06F001 DD DSN=POC3AA7.MRI.MOSESV1.L6(SHUFFLTA),DISP=SHR,
/**          LABEL=(,,,IN)
//FT07F001 DD DUMMY
//FT08F001 DD DUMMY
/**FT08F001 DD DSN=POC3AA7.MRI.MOSESV1.CNTL(DEBUGTB),DISP=SHR,
/**          LABEL=(,,,IN)
//FT69F001 DD DUMMY
/**FT69F001 DD DSN=POC3AA7.MRI.MOSESV1.L69(SHUFFLT1),DISP=SHR,
/**          LABEL=(,,,IN)
//FT09F001 DD SYSOUT=&OUT,DCB=*.FT06F001
//FT90F001 DD DUMMY
//FT99F001 DD SYSOUT=&OUT,DCB=*.FT06F001
/**
//FT47F001 DD DSN=&AEXPO,UNIT=SYSDA,SPACE=(TRK,(40,10)),DISP=(,PASS),

```

```

//          DCB=(RECFM=VBS,LRECL=X,BLKSIZE=2000,BUFNO=2)
//FT41F001 DD DSN=&AEXPR,UNIT=SYSDA,SPACE=(TRK,(100,50)),DISP=(,PASS),
//          DCB=*.FT47F001
//FT49F001 DD DSN=&AEXPR,UNIT=SYSDA,SPACE=(TRK,(40,10)),DISP=(,PASS),
//          DCB=*.FT47F001
//FT48F001 DD DSN=&A6EXP,UNIT=SYSDA,SPACE=(TRK,(40,10)),DISP=(,PASS),
//          DCB=*.FT47F001
//FT72F001 DD DSN=&AEXPR,UNIT=SYSDA,SPACE=(TRK,(40,10)),DISP=(,PASS),
//          DCB=*.FT47F001
//FT73F001 DD DSN=&A6EXR,UNIT=SYSDA,SPACE=(TRK,(40,10)),DISP=(,PASS),
//          DCB=*.FT47F001
/**
//FT12F001 DD DSN=&WKFIL,UNIT=SYSDA,SPACE=(TRK,(5,5)),DISP=(,PASS),
//          DCB=(RECFM=VBS,LRECL=X,BLKSIZE=32760,BUFNO=1)
/**
//          DCB=(RECFM=VBS,LRECL=32756,BLKSIZE=32760,BUFNO=1)
//          DCB=(RECFM=VBS,LRECL=13026,BLKSIZE=13030,BUFNO=1)
//FT13F001 DD DSN=&WKFL0,UNIT=SYSDA,SPACE=(TRK,(5,5)),DISP=(,PASS),
//          DCB=*.FT12F001
/**
//FT20F001 DD DSN=&LOADI,UNIT=SYSDA,SPACE=(TRK,(5,5)),DISP=(,PASS),
//          DCB=*.FT12F001
/**
//FT10F001 DD DSN=&INO66,UNIT=SYSDA,SPACE=(TRK,(40,30)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT11F001 DD DSN=&INO61,UNIT=SYSDA,SPACE=(TRK,(40,30)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT25F001 DD DSN=&IDPEI,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT26F001 DD DSN=&KEFFC,UNIT=SYSDA,SPACE=(TRK,(10,10)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT27F001 DD DSN=&RPEAK,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT28F001 DD DSN=&POFLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT31F001 DD DSN=&AJFLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT32F001 DD DSN=&FTFLY,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT39F001 DD DSN=&P13PO,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT40F001 DD DSN=&P6FLU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT42F001 DD DSN=&DLEAK,UNIT=SYSDA,SPACE=(TRK,(10,10)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT44F001 DD DSN=&AFLUX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT45F001 DD DSN=&AP0WR,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT46F001 DD DSN=&P6POW,UNIT=SYSDA,SPACE=(TRK,(100,25)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT50F001 DD DSN=&AVTEM,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT54F001 DD DSN=&RDEEN,UNIT=SYSDA,SPACE=(TRK,(15,15)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT55F001 DD DSN=&FLUX,UNIT=SYSDA,SPACE=(TRK,(500,500)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT56F001 DD DSN=&PTAU,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=*.FT12F001
//FT57F001 DD DSN=&SOURC,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),

```

```

//          DCB=* .FT12F001
//FT58F001 DD DSN=&PTSAE,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=* .FT12F001
//FT59F001 DD DSN=&DLEK1,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=* .FT12F001
//FT60F001 DD DSN=&FTFLX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=* .FT12F001
//FT61F001 DD DSN=&P6HTX,UNIT=SYSDA,SPACE=(TRK,(20,20)),DISP=(,PASS),
//          DCB=* .FT12F001
//FT16F001 DD DSN=&HIST1,UNIT=SYSDA,SPACE=(TRK,(30,30)),DISP=(,PASS),
//          DCB=* .FT12F001
//FT17F001 DD DSN=&HIST2,UNIT=SYSDA,SPACE=(TRK,(30,30)),DISP=(,PASS),
//          DCB=* .FT12F001
//PLOTLOG DD DUMMY
//GDRILE DD DUMMY
//*PLOTLOG DD SYSOUT=&OUT,DCB=* .FT06F001
//*GDRFILE DD SYSOUT=&OUT,DCB=* .FT06F001
//*PLOTPRM DD *
//* SCALE=0.7
//*-----
//FT01F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=* .FT12F001
//FT02F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT14F001 DD UNIT=SYSDA,SPACE=(TRK,(&N4)),DCB=* .FT01F001
//FT15F001 DD UNIT=SYSDA,SPACE=(TRK,(&N7)),DCB=* .FT01F001
//FT04F001 DD UNIT=SYSDA,SPACE=(TRK,(&N8)),DCB=* .FT15F001
//FT18F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT19F001 DD UNIT=SYSDA,SPACE=(TRK,(&N9)),DCB=* .FT01F001
//FT33F001 DD UNIT=SYSDA,SPACE=(TRK,(&N10)),DCB=* .FT01F001
//FT34F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT35F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=* .FT01F001
//FT36F001 DD UNIT=SYSDA,SPACE=(TRK,(&N11)),DCB=* .FT01F001
//FT37F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=* .FT01F001
//FT38F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT43F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT51F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT52F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,2)),DCB=* .FT01F001
//FT62F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//FT63F001 DD UNIT=SYSDA,SPACE=(TRK,(&NS,1)),DCB=* .FT01F001
//*--- INTERFACE FILE
//FT74F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT75F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT76F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT77F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT79F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT82F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT85F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//FT86F001 DD UNIT=SYSDA,SPACE=(TRK,(&N1,&NX)),DCB=* .FT01F001
//          PEND
//*
//          EXEC MOSESV1,
//          N1=100,N2=6526,N3=10,N4=1,N5=2877,N6=10,
//          N7=103,N8=103,N9=933,N10=2414,N11=1015,N12=1,N13=1,N14=1,
//          N15=10,N16=6526,NX=200,NS=5000
//PLOTLOG DD DUMMY
//GDRFILE DD SYSOUT=R
//PLOTPRM DD *
//          SCALE=0.7
//*
//FT03F001 DD DISP=SHR,DSN=POCOB17.OBHZHBC7.MIC7GJ3,LABEL=(, , IN)

```

```

//FT05F001 DD DISP=SHR,LABEL=(, , IN),
//          DSN=POCOB17.OPMOSD.DATA(GDI1MBRN)
//

```

Example Data Files

PNC TN9460 98-001

Data file [1] .OPDEND.DATA(GDI1P)

PENCIL input, 'iterative condensation' & 'basic fuel cycle'

```
&SET
LAST= 1,
PURICH= 40.2453,45.0307,
COREH= 3.0, 6.0, 9.0, 12.0, 12.0, 9.0, 6.0, 3.0,
ICZ= 1, 2, 3, 4, 5, 6, 7, 8,
OCZ= 9,10,11,12,13,14,15,16,17,18,19,20,21,22,23,24,25,26,27,28,
      29,30,31,32,
NUID= 1, 2, 3, 4, 5, 6, 7, 8, 9, 10,
      11, 12, 13, 14, 15, 0, 0, 18, 0, 20,
      21, 22, 23, 24, 25, 26, 27, 28, 29, 0,
      0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0,
      0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0,
      0, 0, 0, 0, 0, 0, 0, 0, 0, 60,
      0, 0, 0, 0, 0, 0, 67, 68, 0, 70,
      0, 0, 0, 0, 0, 0, 0, 78, 79, 80,
      81, 0, 0, 0, 0,
F=1.0, 1.0, 1.0, 1.0, 1.0, 1.0, 1.0, 1.0,
EPOWER=60.0,
LPOWER=0.0,
RONA=0.837936,
DAY=296.56, 4.0, 6.0,
&END
INNER, CORE 7, HQPU, 4*9.75M., F165%, 24P ZRH, 8P B4C(30%), 45%HOLE, GDI1
&DATA
VB4C=0.060248,
B10W=30.0,
DCLAD=7.422,
HCLAD=0.5376,
HWRAP=3.90,
THDENS=0.687999,
DFUEL=6.141,
DWIRE=0.00,
HWIRE=200.0,
PRING=3.8040,
NOPIN=217,
ITYPE=0,
PUABND=0.94,0.06,0.0,0.0,0.0,0.0,0.0,
UAEND=0.003,0.997,
VSUS=0.6550,0.1700,0.1350,0.0250,0.0150,
ROSUS=7.98,
DGAP=5.8,
SZRH=1.0,
NZRH=24,
ZRH=1.7,
TRRICH=0.0,
TRU=3*0.0,0.491,6*0.0,0.301,0.0,0.155,2*0.0,0.05,0.003,
```

```
&END
OUTER CORE -----
&DATA
VB4C=0.060248,
B10W=30.0,
DCLAD=7.422,
HCLAD=0.5376,
HWRAP=3.90,
THDENS=0.687999,
DFUEL=6.141,
DWIRE=0.00,
HWIRE=200.0,
PRING=3.8040,
NOPIN=217,
ITYPE=0,
PUABND=0.94,0.06,0.0,0.0,0.0,0.0,0.0,
UAEND=0.003,0.997,
VSUS=0.6550,0.1700,0.1350,0.0250,0.0150,
ROSUS=7.98,
DGAP=5.8,
SZRH=1.0,
NZRH=24,
ZRH=1.7,
TRRICH=0.0,
TRU=3*0.0,0.491,6*0.0,0.301,0.0,0.155,2*0.0,0.05,0.003,
&END
&CONV
CPW =1.0,
EPW =0.002,
CKEFF = 1.00324,
EKEFF = 0.0005,
&END
```

Data file [2] .OPCITD.DATA(O975M165)

CITATION input from PENCIL

```
600MW/60CM CORE PU BURNER , OPTIMIZATION , 165% PINSIZE REFLECTORS
296.56*4 DAYS , PENCIL DATA , CORE 7 , FULL HEIGHT & RODS , BCR&PCR 0%
001
1 0 0 0 0 1 1 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
1 0 0 1 0 1 0 1 0 0 0 0 1 1 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
200200
002
6 1 1 1 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
296.6 1560.0
296.6 0.0 0.0 296.6 0.0 0.0
1 0 0 1 0 1 0 1 0 0 0 0 1 1 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
```



```

003
0 0 0 0 7 0 0 0 0 0 1 0 0 0 0 0 0 0 0 0 0 0 0 0
1.0000E-4 1.0000E-5
0.0 0.0 1560.0
004
10 46.21720 1 4.27490 7 38.52480 1 2.29260 1 2.23650 7 38.48690
1 4.61620 6 28.74400 1 1.24510 1 2.46280 5 21.29210 2 6.00000
1 3.00000 4 21.10000 4 21.10000
10 50.00000 2 7.00000 1 3.00000 2 5.00000 1 3.00000 2 6.00000
3 9.00000 4 12.00000 4 12.00000 3 9.00000 2 6.00000 1 3.00000
1 2.00000 14 70.00000 6 33.00000
005
46 49 46 50 50 46 51 46 52 52 46 46 46 46 46
41 49 41 50 50 41 51 41 52 52 41 41 41 47 46
37 49 37 50 50 37 51 38 52 52 38 38 38 47 46
33 49 33 50 50 33 51 34 52 52 34 34 34 47 46
1 42 1 43 43 1 44 9 45 45 9 17 25 47 46
2 42 2 43 43 2 44 10 45 45 10 18 26 47 46
3 42 3 43 43 3 44 11 45 45 11 19 27 47 46
4 42 4 43 43 4 44 12 45 45 12 20 28 47 46
5 42 5 43 43 5 44 13 45 45 13 21 29 47 46
6 42 6 43 43 6 44 14 45 45 14 22 30 47 46
7 42 7 43 43 7 44 15 45 45 15 23 31 47 46
8 42 8 43 43 8 44 16 45 45 16 24 32 47 46
39 42 39 43 43 39 44 40 45 45 40 40 40 47 46
35 42 35 43 43 35 44 36 45 45 36 36 36 47 46
48 48 48 48 48 48 48 48 48 48 48 48 46 46
012
1 1 4 1 0 1 0 INNER CORE1
2 2 4 1 0 2 0 INNER CORE1
3 3 4 1 0 3 0 INNER CORE1
4 4 4 1 0 4 0 INNER CORE1
5 5 4 1 0 5 0 INNER CORE1
6 6 4 1 0 6 0 INNER CORE1
7 7 4 1 0 7 0 INNER CORE1
8 8 4 1 0 8 0 INNER CORE1
9 9 4 2 0 9 0 OUTER CORE1
10 10 4 2 0 10 0 OUTER CORE1
11 11 4 2 0 11 0 OUTER CORE1
12 12 4 2 0 12 0 OUTER CORE1
13 13 4 2 0 13 0 OUTER CORE1
14 14 4 2 0 14 0 OUTER CORE1
15 15 4 2 0 15 0 OUTER CORE1
16 16 4 2 0 16 0 OUTER CORE1
17 17 4 2 0 17 0 OUTER CORE2
18 18 4 2 0 18 0 OUTER CORE2
19 19 4 2 0 19 0 OUTER CORE2
20 20 4 2 0 20 0 OUTER CORE2
21 21 4 2 0 21 0 OUTER CORE2
22 22 4 2 0 22 0 OUTER CORE2
23 23 4 2 0 23 0 OUTER CORE2
24 24 4 2 0 24 0 OUTER CORE2
25 25 4 2 0 25 0 OUTER CORE3
26 26 4 2 0 26 0 OUTER CORE3
27 27 4 2 0 27 0 OUTER CORE3
28 28 4 2 0 28 0 OUTER CORE3
29 29 4 2 0 29 0 OUTER CORE3
30 30 4 2 0 30 0 OUTER CORE3
31 31 4 2 0 31 0 OUTER CORE3

```

```

32 32 4 2 0 32 0 OUTER CORE3
33 33 0 3 0 33 -1 UPPER PLENUM INNER
34 34 0 3 0 34 -1 UPPER PLENUM OUTER
35 35 0 3 0 35 -1 LOWER PLENUM INNER
36 36 0 3 0 36 -1 LOWER PLENUM OUTER
37 37 0 4 0 37 -1 UPPER PLUG INNER
38 38 0 4 0 38 -1 UPPER PLUG OUTER
39 39 0 4 0 39 -1 LOWER PLUG INNER
40 40 0 4 0 40 -1 LOWER PLUG OUTER
41 41 0 5 0 41 -1 UPPER GAP
42 42 0 5 0 42 -1 ROD FOLLOWER 1
43 43 0 5 0 43 -1 ROD FOLLOWER 2
44 44 0 5 0 44 -1 ROD FOLLOWER 3
45 45 0 5 0 45 -1 ROD FOLLOWER 4
46 46 0 6 0 46 -1 UPPER/OUTER SHIELD
47 47 0 6 0 47 -1 RADIAL SHIELD INNER
48 48 0 7 0 48 -1 LOWER SHIELD
49 49 0 8 0 49 -1 ROD ABSORBER 1
50 50 0 8 0 50 -1 ROD ABSORBER 2
51 51 0 8 0 51 -1 ROD ABSORBER 3
52 52 0 8 0 52 -1 ROD ABSORBER 4
018
1 3 0 2 8 0 0 4 5 6 7 21 22 23 24 25 26 27 28 29 67 0 11 12
13 14 15 18 68 0 9 10 0 78 79 80 81 0
020
33 36 0 0
101.03970-2 111.17900-2 123.28650-3 132.31220-3 142.61950-4 152.74470-4
37 40 0 0
101.03970-2 112.96650-2 128.26930-3 135.81800-3 146.59100-4 156.90610-4
41 45 0 0
101.99080-2 115.24170-3 121.46120-3 131.02800-3 141.16460-4 151.22020-4
46 47 0 0
104.38990-3 114.50900-2 121.25690-2 138.84310-3 141.00180-3 151.04970-3
48 48 0 0
106.36540-3 111.77540-2 124.94900-3 133.48200-3 143.94460-4 154.13320-4
189.24600-3 677.89610-3 682.90880-2
49 52 0 0
109.32850-3 111.27380-2 123.55070-3 132.49820-3 142.83010-4 152.96540-4
188.22780-3 673.04990-2 682.41210-3
034
1 2
31.52636-9 223.40600-6 254.92182-8
1 2
1 2 3 4 5 6 7 8 21 22 23 24 25 26 27 28 29
0.0 78 0.0 0.0 0.0 0.0 0.0
0.0 0.0 1.0 0.0 0.0 0.0 0.0
0.0 0.0 0.0 0.0 0.0 0.0 0.0
0.0 79 1.0 0.0 0.0 0.0 0.0
0.0 0.0 0.0 0.0 0.0 0.0 0.0
0.0 0.0 0.0 0.0 0.0 0.0 0.0
0.0 80 0.0 0.0 0.0 0.0 0.0 0.0
1.0 1.0 0.0 0.0 0.0 0.0 0.0

```


Data file [3] .OPEDID.DATA(GDI1RZ)

RZOUT3 input for 'basic fuel cycle'

```

&ZONE
NID= 85,6,9
IDCIT(1,1)= 1, 2, 3, 4, 5, 6, 7, 8, 9, 10,
11, 12, 13, 14, 15, 16, 17, 18, 19, 20,
21, 22, 23, 24, 25, 26, 27, 28, 29, 30,
31, 32, 33, 34, 35, 30, 37, 38, 39, 40,
41, 42, 43, 44, 45, 46, 47, 48, 49, 50,
51, 52, 53, 54, 55, 56, 57, 58, 59, 60,
61, 62, 63, 64, 65, 66, 67, 68, 69, 70,
71, 72, 73, 74, 75, 76, 77, 78, 79, 80,
81, 82, 83, 84, 85, 14*0,
IDSLA(1,1)=949,940,941,942,951,925,926,928, 8, 11,
26, 24, 28, 42, 25,884,894, 6,157, 40,
937,939,950,953,962,963,964,965,948,854,
814,966,967,968,439,579,580,582,584,591,
603,604,605,606,608,600,617,627,628,620,
622,631,633,634,644,646,648, 82,839, 12,
39,441,442,443,529,555,105,115,147, 1,
402, 4, 13, 63, 73, 74, 20,887,897,857,
817,889,899,859,819, 14*0,
IDCIT(1,2)= 10, 11, 12, 13, 14, 15, 93*0,
IDSLA(1,2)= 11, 26, 24, 28, 42, 25, 93*0,
IDCIT(1,3)= 10, 11, 12, 13, 14, 15, 18, 67, 68, 90*0,
IDSLA(1,3)= 11, 26, 24, 28, 42, 25, 6, 105, 115, 90*0,
NEWZN(1)=1,2,3,4,5,6,7,8,9,10,11,12,13,14,15,16,17,18,19,20,21,22,23,
24,25,26,27,28,29,30,31,32,33,34,35,36,37,38,39,40,41,42,43,
44,45,46,47,48,49,50,51,52,38*0,
ZNGR(1)=32*1,15*2,5*3,
ISTEP=17,
NZONE=52, 52,
&END

```

Data file [4] .OPFORT.DATA(GDI1RZ)

output from RZOUT3, for 'basic fuel cycle'

```

POCOB17.@GDI1.DENS
*****CITATION DATA*****
1 1 0 0
11.46918-3 21.74999-4 39.65638-6 43.30258-7 53.79707-7 66.73607-6
73.35554-7 82.60913-3 98.90358-3 101.03968-2 111.17896-2 123.28651-3
132.31220-3 142.61937-4 152.74457-4 160.0 170.0 182.84691-4
190.0 201.28188-3 212.91854-7 223.71285-7 236.62542-9 241.22009-8
251.51758-8 263.8220-10 278.5073-10 282.3774-11 294.90057-8 300.0
310.0 320.0 330.0 340.0 350.0 300.0
370.0 380.0 390.0 400.0 410.0 420.0
430.0 440.0 450.0 460.0 470.0 480.0
490.0 500.0 510.0 520.0 530.0 540.0

```

```

550.0 560.0 570.0 580.0 590.0 600.0
610.0 620.0 630.0 640.0 650.0 660.0
672.83107-4 687.74029-4 690.0 702.17919-3 710.0 720.0
730.0 740.0 750.0 760.0 770.0 789.15262-6
792.09324-4 801.09239-6 815.96565-6 820.0 830.0 840.0
850.0
2 2 0 0
11.44892-3 21.77607-4 31.00720-5 43.65641-7 53.93013-7 66.62858-6
73.51142-7 82.60133-3 98.90358-3 101.03968-2 111.17896-2 123.28651-3
132.31220-3 142.61937-4 152.74457-4 160.0 170.0 182.84691-4
190.0 201.28188-3 213.68157-7 224.00145-7 237.22962-9 241.38310-8
251.66777-8 264.3983-10 271.01782-9 282.9681-11 296.21125-8 300.0
310.0 320.0 330.0 340.0 350.0 300.0
370.0 380.0 390.0 400.0 410.0 420.0
430.0 440.0 450.0 460.0 470.0 480.0
490.0 500.0 510.0 520.0 530.0 540.0
550.0 560.0 570.0 580.0 590.0 600.0
610.0 620.0 630.0 640.0 650.0 660.0
672.78976-4 687.74029-4 690.0 702.17919-3 710.0 720.0
730.0 740.0 750.0 760.0 770.0 781.17643-5
792.30406-4 801.18318-6 817.26002-6 820.0 830.0 840.0
850.0
3 3 0 0
11.41403-3 21.84530-4 31.15877-5 44.78984-7 54.44785-7 66.42750-6
73.88781-7 82.58786-3 98.90358-3 101.03968-2 111.17896-2 123.28651-3

```

REPETITIVE DATA REMOVED

```

730.0 740.0 750.0 760.0 770.0 785.88512-6
791.60877-4 807.03978-7 813.73418-6 820.0 830.0 840.0
850.0
31 31 0 0
11.72898-3 21.72764-4 36.89380-6 41.58220-7 52.83502-7 66.64365-6
72.17290-7 82.42931-3 98.92030-3 101.03968-2 111.17896-2 123.28651-3
132.31221-3 142.61937-4 152.74458-4 160.0 170.0 182.84691-4
190.0 201.28188-3 211.60892-7 222.26524-7 233.37278-9 244.09906-9
257.45128-9 261.2826-10 271.9069-10 283.6340-12 291.98672-8 300.0
310.0 320.0 330.0 340.0 350.0 300.0
370.0 380.0 390.0 400.0 410.0 420.0
430.0 440.0 450.0 460.0 470.0 480.0
490.0 500.0 510.0 520.0 530.0 540.0
550.0 560.0 570.0 580.0 590.0 600.0
610.0 620.0 630.0 640.0 650.0 660.0
673.07470-4 687.74028-4 690.0 702.17919-3 710.0 720.0
730.0 740.0 750.0 760.0 770.0 784.84650-6
791.50914-4 806.71913-7 813.24357-6 820.0 830.0 840.0
850.0
32 32 0 0
11.72436-3 21.77189-4 38.05075-6 42.02331-7 53.28755-7 66.60039-6
72.36147-7 82.42788-3 98.92030-3 101.03968-2 111.17896-2 123.28651-3
132.31221-3 142.61937-4 152.74458-4 160.0 170.0 182.84691-4
190.0 201.28188-3 211.29464-7 222.40311-7 234.26227-9 246.18931-9
259.46179-9 261.8090-10 273.1931-10 286.8851-12 292.01640-8 300.0
310.0 320.0 330.0 340.0 350.0 300.0
370.0 380.0 390.0 400.0 410.0 420.0

```

430.0	440.0	450.0	460.0	470.0	480.0
490.0	500.0	510.0	520.0	530.0	540.0
550.0	560.0	570.0	580.0	590.0	600.0
610.0	620.0	630.0	640.0	650.0	660.0
673.02455-4	687.74029-4	690.0	702.17919-3	710.0	720.0
730.0	740.0	750.0	760.0	770.0	783.83541-6
791.52442-4	806.94562-7	813.06902-6	820.0	830.0	840.0
850.0					
33 33 0 0					
101.03970-2	111.17900-2	123.28650-3	132.31220-3	142.61950-4	152.74470-4
34 34 0 0					
101.03970-2	111.17900-2	123.28650-3	132.31220-3	142.61950-4	152.74470-4
35 35 0 0					
101.03970-2	111.17900-2	123.28650-3	132.31220-3	142.61950-4	152.74470-4
36 36 0 0					
101.03970-2	111.17900-2	123.28650-3	132.31220-3	142.61950-4	152.74470-4
37 37 0 0					
101.03970-2	112.96650-2	128.26930-3	135.81799-3	146.59100-4	156.90610-4
38 38 0 0					
101.03970-2	112.96650-2	128.26930-3	135.81799-3	146.59100-4	156.90610-4
39 39 0 0					
101.03970-2	112.96650-2	128.26930-3	135.81799-3	146.59100-4	156.90610-4
40 40 0 0					
101.03970-2	112.96650-2	128.26930-3	135.81799-3	146.59100-4	156.90610-4
41 41 0 0					
101.99080-2	115.24170-3	121.46120-3	131.02800-3	141.16460-4	151.22020-4
42 42 0 0					
101.99080-2	115.24170-3	121.46120-3	131.02800-3	141.16460-4	151.22020-4
43 43 0 0					
101.99080-2	115.24170-3	121.46120-3	131.02800-3	141.16460-4	151.22020-4
44 44 0 0					
101.99080-2	115.24170-3	121.46120-3	131.02800-3	141.16460-4	151.22020-4
45 45 0 0					
101.99080-2	115.24170-3	121.46120-3	131.02800-3	141.16460-4	151.22020-4
46 46 0 0					
104.38989-3	114.50900-2	121.25690-2	138.84309-3	141.00180-3	151.04970-3
47 47 0 0					
104.38990-3	114.50900-2	121.25690-2	138.84310-3	141.00180-3	151.04970-3
48 48 0 0					
106.36540-3	111.77540-2	124.94900-3	133.48200-3	143.94460-4	154.13320-4
189.24600-3	677.89610-3	682.90880-2			
49 49 0 0					
109.32850-3	111.27380-2	123.55070-3	132.49820-3	142.83010-4	152.96540-4
188.22780-3	673.04990-2	682.41210-3			

```

50 50 0 0
109.32850-3 111.27380-2 123.55070-3 132.49820-3 142.83010-4 152.96540-4
188.22780-3 673.04990-2 682.41210-3
51 51 0 0
109.32850-3 111.27380-2 123.55070-3 132.49820-3 142.83010-4 152.96540-4
188.22780-3 673.04990-2 682.41210-3
52 52 0 0
109.32850-3 111.27380-2 123.55070-3 132.49820-3 142.83010-4 152.96540-4
188.22780-3 673.04990-2 682.41210-3
*****END*****
*****SLAROM DATA*****

```

```

ZONE 1
949 1.46918E-03 940 1.74999E-04 941 9.65638E-06 942 3.30258E-07
951 3.79707E-07 925 6.73607E-06 926 3.35554E-07 928 2.60913E-03
8 8.90358E-03 11 1.03968E-02 26 1.17896E-02 24 3.28651E-03
28 2.31220E-03 42 2.61937E-04 25 2.74457E-04 884 0.00000E+00
894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28188E-03
937 2.91854E-07 939 3.71285E-07 950 6.62542E-09 953 1.22009E-08
962 1.51758E-08 963 3.82196E-10 964 8.50733E-10 965 2.37739E-11
948 4.90057E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
529 0.00000E+00 555 0.00000E+00 105 2.83107E-04 115 7.74029E-04
147 0.00000E+00 1 2.17919E-03 402 0.00000E+00 4 0.00000E+00
13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
20 0.00000E+00 887 9.15262E-06 897 2.09324E-04 857 1.09239E-06
817 5.96565E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
819 0.00000E+00

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ZONE 2
949 1.44892E-03 940 1.77607E-04 941 1.00720E-05 942 3.65641E-07
951 3.93013E-07 925 6.62858E-06 926 3.51142E-07 928 2.60133E-03
8 8.90358E-03 11 1.03968E-02 26 1.17896E-02 24 3.28651E-03
28 2.31220E-03 42 2.61937E-04 25 2.74457E-04 884 0.00000E+00
894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28188E-03
937 3.68157E-07 939 4.00145E-07 950 7.22962E-09 953 1.38310E-08
962 1.66777E-08 963 4.39829E-10 964 1.01782E-09 965 2.96814E-11
948 6.21125E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
529 0.00000E+00 555 0.00000E+00 105 2.78976E-04 115 7.74029E-04
147 0.00000E+00 1 2.17919E-03 402 0.00000E+00 4 0.00000E+00
13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
20 0.00000E+00 887 1.17643E-05 897 2.30406E-04 857 1.18318E-06
817 7.26002E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
819 0.00000E+00

```

ZONE 3
949 1.41403E-03 940 1.84530E-04 941 1.15877E-05 942 4.78984E-07
951 4.44785E-07 925 6.42750E-06 926 3.88781E-07 928 2.58786E-03

.....
REPETITIVE DATA REMOVED
.....

20 0.00000E+00 887 4.84650E-06 897 1.50914E-04 857 6.71913E-07
817 3.24357E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
819 0.00000E+00

ZONE 32
949 1.72436E-03 940 1.77189E-04 941 8.05075E-06 942 2.02331E-07
951 3.28755E-07 925 6.60039E-06 926 2.36147E-07 928 2.42788E-03
8 8.92030E-03 11 1.03968E-02 26 1.17896E-02 24 3.28651E-03
28 2.31221E-03 42 2.61937E-04 25 2.74458E-04 884 0.00000E+00
894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28188E-03
937 1.29464E-07 939 2.40311E-07 950 4.26227E-09 953 6.18931E-09
962 9.46179E-09 963 1.80899E-10 964 3.19306E-10 965 6.88508E-12
948 2.01640E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
529 0.00000E+00 555 0.00000E+00 105 3.02455E-04 115 7.74029E-04
147 0.00000E+00 1 2.17919E-03 402 0.00000E+00 4 0.00000E+00
13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
20 0.00000E+00 887 3.83541E-06 897 1.52442E-04 857 6.94562E-07
817 3.06902E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
819 0.00000E+00

ZONE 33
11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
42 2.61950E-04 25 2.74470E-04

ZONE 34
11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
42 2.61950E-04 25 2.74470E-04

ZONE 35

.....
REPETITIVE DATA REMOVED
.....

ZONE 47
11 4.38990E-03 26 4.50900E-02 24 1.25690E-02 28 8.84310E-03
42 1.00180E-03 25 1.04970E-03

ZONE 48
11 6.36540E-03 26 1.77540E-02 24 4.94900E-03 28 3.48200E-03
42 3.94460E-04 25 4.13320E-04 6 9.24600E-03 105 7.89610E-03
115 2.90880E-02

ZONE 49
11 9.32850E-03 26 1.27380E-02 24 3.55070E-03 28 2.49820E-03
42 2.83010E-04 25 2.96540E-04 6 8.22780E-03 105 3.04990E-02
115 2.41210E-03

ZONE 50
11 9.32850E-03 26 1.27380E-02 24 3.55070E-03 28 2.49820E-03
42 2.83010E-04 25 2.96540E-04 6 8.22780E-03 105 3.04990E-02
115 2.41210E-03

ZONE 51
11 9.32850E-03 26 1.27380E-02 24 3.55070E-03 28 2.49820E-03
42 2.83010E-04 25 2.96540E-04 6 8.22780E-03 105 3.04990E-02
115 2.41210E-03

ZONE 52
11 9.32850E-03 26 1.27380E-02 24 3.55070E-03 28 2.49820E-03
42 2.83010E-04 25 2.96540E-04 6 8.22780E-03 105 3.04990E-02
115 2.41210E-03

*****END*****

Data file [5] .OPDEND.DATA(GDG0S)

SLAROM input for 'iterative condensation'

PREP

INNER CORE 1,CORE 7,HQPU39.86/44.86%,4*9.75,165%,31/42RH,BOC GDGO
2 1 0 1 1 0 3 -20 0 0 0 0 70
10703.15 1.3 0.0
1373.15 703.15

77 8
0.30705 0.30160
949 0.549351E-02 940 0.735227E-03 941 0.486471E-04 942 0.214555E-05
951 0.185357E-05 925 0.252666E-04 926 0.161152E-05 928 0.102704E-01
8 0.351686E-01 884 0.000000E+00 894 0.000000E+00 6 0.111864E-02
157 0.000000E+00 937 0.179786E-05 939 0.191532E-05 950 0.406384E-07
953 0.984156E-07 962 0.959287E-07 963 0.304819E-08 964 0.895688E-08
965 0.322166E-09 948 0.363717E-06 854 0.000000E+00 814 0.000000E+00
966 0.000000E+00 967 0.000000E+00 968 0.000000E+00 439 0.000000E+00
579 0.000000E+00 580 0.000000E+00 582 0.000000E+00 584 0.000000E+00
591 0.000000E+00 603 0.000000E+00 604 0.000000E+00 605 0.000000E+00
606 0.000000E+00 608 0.000000E+00 600 0.000000E+00 617 0.000000E+00
627 0.000000E+00 628 0.000000E+00 620 0.000000E+00 622 0.000000E+00
631 0.000000E+00 633 0.000000E+00 634 0.000000E+00 644 0.000000E+00
646 0.000000E+00 648 0.000000E+00 82 0.000000E+00 839 0.000000E+00
12 0.000000E+00 39 0.000000E+00 441 0.000000E+00 442 0.000000E+00
443 0.000000E+00 529 0.000000E+00 555 0.000000E+00 105 0.104222E-02
115 0.304139E-02 147 0.000000E+00 402 0.000000E+00 4 0.000000E+00
13 0.000000E+00 63 0.000000E+00 73 0.000000E+00 74 0.000000E+00

20 0.000000E+00 887 0.591615E-04 897 0.107694E-02 857 0.554935E-05
817 0.393071E-04 889 0.000000E+00 899 0.000000E+00 859 0.000000E+00
819 0.000000E+00

11 0.139463E-01 26 0.158146E-01 24 0.440851E-02 28 0.310159E-02
42 0.351362E-03 25 0.368157E-03 40 0.164787E-02 1 0.280137E-02

PATH

HETEROGENEOUS 2 TEMPERATURE MODEL

70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0

1 2
0.0 0.30705 0.60865

PIJF

37 0 5 3 0 0 0 0
0 0 0. 0. 0. 0.

EDIT

2 2 1 0 0

INCORE

PREP

OUTER CORE 2,CORE 7,HQPU39.86/44.86%,4*9.75,165%,31/4ZRH,BOC GDGO

2 1 0 1 1 0 3 -20 0 0 0 0 70

10703.15 1.3 0.0

1373.15 703.15

77 8
0.30705 0.26171

949 0.567654E-02 940 0.637067E-03 941 0.306210E-04 942 0.958221E-06

951 0.122012E-05 925 0.218034E-04 926 0.957798E-06 928 0.833109E-02

8 0.307707E-01 884 0.000000E+00 894 0.000000E+00 6 0.976832E-03

157 0.000000E+00 937 0.115092E-05 939 0.109166E-05 950 0.193127E-07

953 0.306155E-07 962 0.439324E-07 963 0.100419E-08 964 0.197616E-08

965 0.510089E-10 948 0.165980E-06 854 0.000000E+00 814 0.000000E+00

966 0.000000E+00 967 0.000000E+00 968 0.000000E+00 439 0.000000E+00

579 0.000000E+00 580 0.000000E+00 582 0.000000E+00 584 0.000000E+00

591 0.000000E+00 603 0.000000E+00 604 0.000000E+00 605 0.000000E+00

606 0.000000E+00 608 0.000000E+00 600 0.000000E+00 617 0.000000E+00

627 0.000000E+00 628 0.000000E+00 620 0.000000E+00 622 0.000000E+00

631 0.000000E+00 633 0.000000E+00 634 0.000000E+00 644 0.000000E+00

646 0.000000E+00 648 0.000000E+00 82 0.000000E+00 839 0.000000E+00

12 0.000000E+00 39 0.000000E+00 441 0.000000E+00 442 0.000000E+00

443 0.000000E+00 529 0.000000E+00 555 0.000000E+00 105 0.100026E-02

115 0.265585E-02 147 0.000000E+00 402 0.000000E+00 4 0.000000E+00

13 0.000000E+00 63 0.000000E+00 73 0.000000E+00 74 0.000000E+00

20 0.000000E+00 887 0.365056E-04 897 0.775911E-03 857 0.329013E-05

817 0.236091E-04 889 0.000000E+00 899 0.000000E+00 859 0.000000E+00

819 0.000000E+00

11 0.146730E-01 26 0.166387E-01 24 0.463825E-02 28 0.326322E-02

42 0.369672E-03 25 0.387342E-03 40 0.173374E-02 1 0.294733E-02

PATH

HETEROGENEOUS 2 TEMPERATURE MODEL

70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0

1 2
0.0 0.30705 0.56876

PIJF

37 0 5 3 0 0 0 0

0 0 0. 0. 0. 0.

EDIT

2 2 1 0 0

OTCORE

PREP

PLENUM UP&LO, LO Q PU,165%FUEL,

1 1 0 0 0 0 3 -20 0 0 0 0 70

703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03

42 2.61950E-04 25 2.74470E-04

PLENUM

PREP

PLUGGED PIN, LO Q ,165%FUEL

1 1 0 0 0 0 3 -20 0 0 0 0 70

703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 2.96650E-02 24 8.26930E-03 28 5.81800E-03

42 6.59100E-04 25 6.90610E-04

PLUGGD

PREP

ROD FOLLOWER/GAP,

1 1 0 0 0 0 3 -20 0 0 0 0 70

703.15 0.0 0.0

6

1.0

11 1.99080E-02 26 5.24170E-03 24 1.46120E-03 28 1.02800E-03

42 1.16460E-04 25 1.22020E-04

FOLLOW

PREP

UPP/RAD'L SH'LD,

1 1 0 0 0 0 3 -20 0 0 0 0 70

703.15 0.0 0.0

6

1.0

11 4.38990E-03 26 4.50900E-02 24 1.25690E-02 28 8.84310E-03

42 1.00180E-03 25 1.04970E-03

UPPSHD

PREP

LOWER SHIELD,

1 1 0 0 0 0 3 -20 0 0 0 0 70

703.15 0.0 0.0

9

1.0

11 6.36540E-03 26 1.77540E-02 24 4.94900E-03 28 3.48200E-03

42 3.94460E-04 25 4.13320E-04 6 9.24600E-03 105 7.89610E-03

115 2.90880E-02

LOWSHD

PREP

ABSORBER ROD, HETEROGENEOUS CELL MODEL WITHIN 9.7 S/A FUEL

6 1 0 1 1 0 3 -21 0 0 0 0 70

703.15 1.3 0.0

3 1 5 1 5 30

4.7901 0.1137 0.9751 2.1419 0.3937 18.8519

6 2.47080E-02 105 9.15890E-02 115 7.24350E-03 /

11 2.19495E-09 /

INCOR
 OTCOR
 PLENU
 PLUGG
 FOLLO
 UPPSH
 LOWSH
 ABSOR

00000120
 00000120
 00000120
 00000120
 00000120
 00000120
 00000120
 00000120
 00000120

ROD ABS-HOMO OPTIMIZED CORE

00000020

PNC TN9460 98-001

Data file [8] .OPEDID.DATA(PFC1C)

condensation JOINT input, 7 groups, for
'iterative condensation'

CITATIONPOST	00000010
1	00000020
7	00000030
4 8 19 28 37 46 70	00000040
PDSIN INCOR SLAR 0 1	00000050
PDSOUT INCOR SLAR	00000060
USERPDS INCOR CITA	00000070
INNER-CORE OPTIMIZED CORE	00000080
PDSIN OTCOR SLAR 0 1	00000020
PDSOUT OTCOR SLAR	00000060
USERPDS OTCOR CITA	00000070
OUTER-CORE OPTIMIZED CORE	00000080
PDSIN PLENU SLAR 0 1	00000020
PDSOUT PLENU SLAR	00000060
USERPDS PLENU CITA	00000070
PLENUM OPTIMIZED CORE	00000080
PDSIN PLUGG SLAR 0 1	00000020
PDSOUT PLUGG SLAR	00000060
USERPDS PLUGG CITA	00000070
STEEL PLUG OPTIMIZED CORE	00000080
PDSIN FOLLO SLAR 0 1	00000020
PDSOUT FOLLO SLAR	00000060
USERPDS FOLLO CITA	00000070
ROD FOLLOWER OPTIMIZED CORE	00000080
PDSIN UPPSH SLAR 0 1	00000020
PDSOUT UPPSH SLAR	00000060
USERPDS UPPSH CITA	00000070
UPPER-SH'D OPTIMIZED CORE	00000080
PDSIN LOWSH SLAR 0 1	00000020
PDSOUT LOWSH SLAR	00000060
USERPDS LOWSH CITA	00000070
RAD'L-SH'D OPTIMIZED CORE	00000080
PDSIN ABSOR SLAR 0 1	00000020
PDSOUT ABSOR SLAR	00000060
USERPDS ABSOR CITA	00000070
ROD ABS-HET OPTIMIZED CORE	00000080
PDSIN AHOMO SLAR 0 1	00000020
PDSOUT AHOMO SLAR	00000060
USERPDS ABSOR CITA	00000070

Data file [9] .OPEDID.DATA(PFC1M)

x-section re-formatting JOINT input for
'iterative condensation'

CITATIONMICRO	00000100
INCOR SLAROM	00000200
85	00000300
1 949 PU239	239.0530 0.334000E-10
2 940 PU240	240.0540 0.336 E-10
3 941 PU241	241.0570 0.337 E-10
4 942 PU242	242.0580 0.338 E-10
5 951 AM241	241.0750 0.336 E-10
6 925 U235	235.0440 0.323 E-10
7 926 U236	236.0460 0.324 E-10
8 928 U238	238.0510 0.331 E-10
9 8 O	15.9994 0.0
10 11 NA	22.9898 0.0
11 26 FE	55.8470 0.0
12 24 CR	51.9961 0.0
13 28 NI	58.6900 0.0
14 42 MO	95.9400 0.0
15 25 MN	54.9380 0.0
16 884 U-238FP	238.0 0.0
17 894 PU-239FP	239.0 0.0
18 6 C	12.011 0.0
19 157 N15	15.000 0.0
20 40 ZR	91.224 0.0
21 937 NP237	237.048 0.330000E-10
22 939 NP239	239.053 0.331000E-10
23 950 AM242M	242.060 0.345000E-10
24 953 AM243	243.061 0.346000E-10
25 962 CM242	242.059 0.352000E-10
26 963 CM243	243.060 0.352000E-10
27 964 CM244	244.063 0.353000E-10
28 965 CM245	245.066 0.353000E-10
29 948 PU238	238.050 0.337000E-10
30 854 U235-FP	235.0 0.0
31 814 PU241-FP	241.0 0.0
32 966 CM246	246.0 0.354000E-10
33 967 CM247	247.0 0.354000E-10
34 968 CM248	248.0 0.354000E-10
35 439 TC99	99.0 0.0
36 579 LA139	139.0 0.0
37 580 CE140	140.0 0.0
38 582 CE142	142.0 0.0
39 584 CE144	144.0 0.0
40 591 PR141	141.0 0.0
41 603 ND143	143.0 0.0
42 604 ND144	144.0 0.0
43 605 ND145	145.0 0.0
44 606 ND146	146.0 0.0

45	608	ND148	148.0	0.0
46	600	ND150	150.0	0.0
47	617	PM147	146.915	0.0
48	627	SM147	147.0	0.0
49	628	SM148	148.0	0.0
50	620	SM150	150.0	0.0
51	622	SM152	152.0	0.0
52	631	EUI151	150.919	0.0
53	633	EUI153	152.921	0.0
54	634	EUI154	154.0	0.0
55	644	GD154	154.0	0.0
56	646	GD156	155.922	0.0
57	648	GD158	157.924	0.0
58	82	PB56	207.2	0.0
59	839	BI209	209.0	0.0
60	12	MG	24.305	0.0
61	39	Y	88.9069	0.0
62	441	RUI01	100.906	0.0
63	442	RUI02	101.904	0.0
64	443	RUI03	102.906	0.0
65	529	I129	128.905	0.0
66	555	CSI35	134.906	0.0
67	105	B10	10.0129	0.0
68	115	B11	11.0093	0.0
69	147	NI4	14.0031	0.0
70	1	H	1.0078	0.0
71	402	HE	4.0026	0.0
72	4	BE	9.0122	0.0
73	13	AL27	26.9818	0.0
74	63	LI6	6.0151	0.0
75	73	LI7	7.0160	0.0
76	74	W-NAT	183.8500	0.0
77	20	CA	40.0800	0.0
78	887	U238-FP	238.0	0.0
79	897	FU239-FP	239.0	0.0
80	857	U235-FP	235.0	0.0
81	817	FU241-FP	241.0	0.0
82	889	U238-FP	238.0	0.0
83	899	FU239-FP	239.0	0.0
84	859	U235-FP	235.0	0.0
85	819	FU241-FP	241.0	0.0

00000420
00000420

INCR 85 INNE CORE REGION
 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 2400000280
 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48
 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72
 73 74 75 76 77 78 79 80 81 82 83 84 85

OTCOR 85 OUT CORE REGION
 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 2400000280
 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44 45 46 47 48
 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 67 68 69 70 71 72
 73 74 75 76 77 78 79 80 81 82 83 84 85

PLENU 6 PLENUM 00000310
 10 11 12 13 14 15 00000280
 PLUGS 6 STEEL PLUG 00000310
 10 11 12 13 14 15 00000280
 FOLLO 6 ROD FOLLOWER 00000320
 10 11 12 13 14 15 00000280
 UPPSH 6 UPPER SHIELD 00000310
 10 11 12 13 14 15 00000280

LOWSH 9 LOWER SHIELD
 10 11 12 13 14 15 18 67 68
 ABSOR 9 ROD ABS-HET
 10 11 12 13 14 15 18 67 68
 AHOMO 9 ROD ABS-HOMO
 10 11 12 13 14 15 18 67 68

00000310
00000280
00000320
00000280
00000320
00000280
00000900

Data file [10] .OPDEND.DATA(GDI1S)

SLAROM input for 'basic fuel cycle'

PREP
 INNER CORE 1,C7,HPU40.25/45.03%,4*9.75M,165%,24/8ZRH/B4CP,SOC GDI1
 2 1 0 1 1 0 3 0 0 0 0 0 70
 10703.15 1.3 0.0
 1373.15 703.15
 77 8
 0.30705 0.30160
 949 0.577285E-02 940 0.687624E-03 941 0.379428E-04 942 0.129768E-05
 951 0.149198E-05 925 0.264680E-04 926 0.131849E-05 928 0.102521E-01
 8 0.349848E-01 884 0.000000E+00 894 0.000000E+00 6 0.111864E-02
 157 0.000000E+00 937 0.114678E-05 939 0.145889E-05 950 0.260333E-07
 953 0.479410E-07 962 0.596302E-07 963 0.150176E-08 964 0.334279E-08
 965 0.934148E-10 948 0.192558E-06 854 0.000000E+00 814 0.000000E+00
 966 0.000000E+00 967 0.000000E+00 968 0.000000E+00 439 0.000000E+00
 579 0.000000E+00 580 0.000000E+00 582 0.000000E+00 584 0.000000E+00
 591 0.000000E+00 603 0.000000E+00 604 0.000000E+00 605 0.000000E+00
 606 0.000000E+00 608 0.000000E+00 600 0.000000E+00 617 0.000000E+00
 627 0.000000E+00 628 0.000000E+00 620 0.000000E+00 622 0.000000E+00
 631 0.000000E+00 633 0.000000E+00 634 0.000000E+00 644 0.000000E+00
 646 0.000000E+00 648 0.000000E+00 82 0.000000E+00 839 0.000000E+00
 12 0.000000E+00 39 0.000000E+00 441 0.000000E+00 442 0.000000E+00
 443 0.000000E+00 529 0.000000E+00 555 0.000000E+00 105 0.111241E-02
 115 0.304139E-02 147 0.000000E+00 402 0.000000E+00 4 0.000000E+00
 13 0.000000E+00 63 0.000000E+00 73 0.000000E+00 74 0.000000E+00
 20 0.000000E+00 887 0.359634E-04 897 0.822497E-03 857 0.429233E-05
 817 0.234408E-04 889 0.000000E+00 899 0.000000E+00 859 0.000000E+00
 819 0.000000E+00

11 0.139463E-01 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
 42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

PATH
 HETEROGENEOUS 2 TEMPERATURE MODEL
 70 2 2 1 0 0 0
 3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
 1 2
 0.0 0.30705 0.60865
 PIJF
 37 0 5 3 0 0 0 0
 0 0 0. 0. 0. 0.
 0
 EDIT

2 2 1 0 0
IC01FF
PREP
INNER CORE 2,C7,HPU40.25/45.03%,4*9.75M,165%,24/8ZRH/B4CP,SOC GDII
2 1 0 1 1 0 3 0 0 0 0 0 70

10703.15 1.3 0.0
1373.15 703.15
77 8
0.30705 0.30160
949 0.569324E-02 940 0.697871E-03 941 0.395759E-04 942 0.143671E-05
951 0.154427E-05 925 0.260457E-04 926 0.137974E-05 928 0.102214E-01
8 0.349848E-01 884 0.000000E+00 894 0.000000E+00 6 0.111864E-02
157 0.000000E+00 937 0.144660E-05 939 0.157229E-05 950 0.284073E-07
953 0.543461E-07 962 0.655317E-07 963 0.172822E-08 964 0.399932E-08
965 0.116627E-09 948 0.244059E-06 854 0.000000E+00 814 0.000000E+00
966 0.000000E+00 967 0.000000E+00 968 0.000000E+00 439 0.000000E+00
579 0.000000E+00 580 0.000000E+00 582 0.000000E+00 584 0.000000E+00
591 0.000000E+00 603 0.000000E+00 604 0.000000E+00 605 0.000000E+00
606 0.000000E+00 608 0.000000E+00 600 0.000000E+00 617 0.000000E+00
627 0.000000E+00 628 0.000000E+00 620 0.000000E+00 622 0.000000E+00
631 0.000000E+00 633 0.000000E+00 634 0.000000E+00 644 0.000000E+00
646 0.000000E+00 648 0.000000E+00 82 0.000000E+00 839 0.000000E+00
12 0.000000E+00 39 0.000000E+00 441 0.000000E+00 442 0.000000E+00
443 0.000000E+00 529 0.000000E+00 555 0.000000E+00 105 0.109618E-02
115 0.304139E-02 147 0.000000E+00 402 0.000000E+00 4 0.000000E+00
13 0.000000E+00 63 0.000000E+00 73 0.000000E+00 74 0.000000E+00
20 0.000000E+00 887 0.462255E-04 897 0.905334E-03 857 0.464907E-05
817 0.285268E-04 889 0.000000E+00 899 0.000000E+00 859 0.000000E+00
819 0.000000E+00
11 0.139463E-01 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

PATH
HETEROGENEOUS 2 TEMPERATURE MODEL
70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
1 2
0.0 0.30705 0.60865
PIJF
37 0 5 3 0 0 0 0
0 0 0. 0. 0. 0.
0
EDIT
2 2 1 0 0
IC02FF
PREP
INNER CORE 3,C7,HPU40.25/45.03%,4*9.75M,165%,24/8ZRH/B4CP,SOC GDII
2 1 0 1 1 0 3 0 0 0 0 0 70

.....
REPETITIVE DATA REMOVED
.....

OC31FF
PREP
OUTER CORE 32,C7,HPU40.25/45.03%,4*9.75M,165%,24/8ZRH/B4CP,SOC GDII
2 1 0 1 1 0 3 0 0 0 0 0 70

10703.15 1.3 0.0
1373.15 703.15
77 8
0.30705 0.26171
949 0.591662E-02 940 0.607971E-03 941 0.276237E-04 942 0.694238E-06
951 0.112802E-05 925 0.226472E-04 926 0.810268E-06 928 0.833054E-02
8 0.306073E-01 884 0.000000E+00 894 0.000000E+00 6 0.976832E-03
157 0.000000E+00 937 0.444217E-06 939 0.824555E-06 950 0.146247E-07
953 0.212368E-07 962 0.324653E-07 963 0.620701E-09 964 0.109560E-08
965 0.236241E-10 948 0.691867E-07 854 0.000000E+00 814 0.000000E+00
966 0.000000E+00 967 0.000000E+00 968 0.000000E+00 439 0.000000E+00
579 0.000000E+00 580 0.000000E+00 582 0.000000E+00 584 0.000000E+00
591 0.000000E+00 603 0.000000E+00 604 0.000000E+00 605 0.000000E+00
606 0.000000E+00 608 0.000000E+00 600 0.000000E+00 617 0.000000E+00
627 0.000000E+00 628 0.000000E+00 620 0.000000E+00 622 0.000000E+00
631 0.000000E+00 633 0.000000E+00 634 0.000000E+00 644 0.000000E+00
646 0.000000E+00 648 0.000000E+00 82 0.000000E+00 839 0.000000E+00
12 0.000000E+00 39 0.000000E+00 441 0.000000E+00 442 0.000000E+00
443 0.000000E+00 529 0.000000E+00 555 0.000000E+00 105 0.103778E-02
115 0.265585E-02 147 0.000000E+00 402 0.000000E+00 4 0.000000E+00
13 0.000000E+00 63 0.000000E+00 73 0.000000E+00 74 0.000000E+00
20 0.000000E+00 887 0.131601E-04 897 0.523059E-03 857 0.238318E-05
817 0.105304E-04 889 0.000000E+00 899 0.000000E+00 859 0.000000E+00
819 0.000000E+00

11 0.146730E-01 26 0.166387E-01 24 0.463825E-02 28 0.326322E-02
42 0.369672E-03 25 0.387342E-03 40 0.180912E-02 1 0.307549E-02

PATH
HETEROGENEOUS 2 TEMPERATURE MODEL
70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
1 2
0.0 0.30705 0.56876
PIJF
37 0 5 3 0 0 0 0
0 0 0. 0. 0. 0.
0

EDIT
2 2 1 0 0
QC32FF
PREP
PLENUM ,LO Q PU,165%FUEL, UPPER INNER
1 1 0 0 0 0 3 0 0 0 0 0 70
703.15 0.0 0.0
6
1.0
11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
42 2.61950E-04 25 2.74470E-04

UIPLNN
PREP
PLENUM ,LO Q PU,165%FUEL, UPPER OUTER
1 1 0 0 0 0 3 0 0 0 0 0 70
703.15 0.0 0.0
6
1.0
11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
42 2.61950E-04 25 2.74470E-04
UOPLNN

PREP

REPETITIVE DATA REMOVED

IRDSHD

PREP

LOWER SHIELD,

1 1 0 0 0 0 3 0 0 0 0 0 70

703.15 0.0 0.0

9

1.0

11 6.36540E-03 26 1.77540E-02 24 4.94900E-03 28 3.48200E-03

42 3.94460E-04 25 4.13320E-04 6 9.24600E-03 105 7.89610E-03

115 2.90880E-02

LSHLDD

PREP

ABSORBER ROD 1, HETEROGENEOUS CELL MODEL WITHIN 9.7 S/A FUEL

6 1 0 1 1 0 3 0 0 0 0 0 70

703.15 1.3 0.0

3 1 5 1 5 30

4.7901 0.1137 0.9751 2.1419 0.3937 18.8519

6 2.47080E-02 105 9.15890E-02 115 7.24350E-03 /

11 2.19495E-09 /

42 1.25230E-03 25 1.31210E-03 /

11 2.19495E-02 /

26 5.63630E-02 24 1.57110E-02 28 1.10540E-02

42 1.25230E-03 25 1.31210E-03 /

949 6.49311E-04 940 8.17696E-04 941 1.68546E-04 942 3.13512E-04

951 1.94612E-05 925 7.78987E-06 926 3.61466E-07 928 2.99340E-03

8 1.03332E-02 11 1.18383E-02 26 1.07946E-02 24 3.00913E-03

28 2.11706E-03 42 2.39830E-04 25 2.51294E-04

40 6.79872E-04

937 3.26805E-07 939 6.64538E-07 950 4.94608E-07 953 1.17133E-05

962 1.13309E-06 963 4.51328E-08 964 1.29052E-06 965 4.89657E-08

948 6.45035E-05

1 1.15578E-03

887 1.02518E-05 897 8.98626E-05 857 1.21029E-06

817 6.73348E-05 /

PATH

TITLE CARD HETEROGENEOUS ROD

70 6 2 1 1 0 0

3 6 6 0 6 0 0 0 0 8 10 0 0 15 0

1 2 3 4 5 6

0.0 4.7901 4.9038 5.7652 7.9071 8.3008 27.1527

PIJF

37 0 5 3 0 1 0 0

0 0 0. 0. 0. 0.

0

EDIT

2 2 1 1 0

1 5 1 5

RODI1A

PREP

ABSORBER ROD 2, HETEROGENEOUS CELL MODEL WITHIN 9.7 S/A FUEL

6 1 0 1 1 0 3 0 0 0 0 0 70

703.15 1.3 0.0

3 1 5 1 5 30

4.7901 0.1137 0.9751 2.1419 0.3937 18.8519

6 2.47080E-02 105 9.15890E-02 115 7.24350E-03 /

11 2.19495E-09 /

42 1.25230E-03 25 1.31210E-03 /

11 2.19495E-02 /

26 5.63630E-02 24 1.57110E-02 28 1.10540E-02

42 1.25230E-03 25 1.31210E-03 /

949 6.49311E-04 940 8.17696E-04 941 1.68546E-04 942 3.13512E-04

951 1.94612E-05 925 7.78987E-06 926 3.61466E-07 928 2.99340E-03

8 1.03332E-02 11 1.18383E-02 26 1.07946E-02 24 3.00913E-03

28 2.11706E-03 42 2.39830E-04 25 2.51294E-04

40 6.79872E-04

937 3.26805E-07 939 6.64538E-07 950 4.94608E-07 953 1.17133E-05

962 1.13309E-06 963 4.51328E-08 964 1.29052E-06 965 4.89657E-08

948 6.45035E-05

1 1.15578E-03

887 1.02518E-05 897 8.98626E-05 857 1.21029E-06

817 6.73348E-05 /

PATH

TITLE CARD HETEROGENEOUS ROD

70 6 2 1 1 0 0

3 6 6 0 6 0 0 0 0 8 10 0 0 15 0

1 2 3 4 5 6

0.0 4.7901 4.9038 5.7652 7.9071 8.3008 27.1527

PIJF

37 0 5 3 0 1 0 0

0 0 0. 0. 0. 0.

0

EDIT

2 2 1 1 0

1 5 1 5

RODI2A

PREP

REPETITIVE DATA REMOVED

RODI4A

Data file [11] .OPCITD.DATA(PEE1D700)

CITATION input for 'basic fuel cycle'

CITATION IC01FF SLAROM

00000010 00000120


```

42
FOLL1 SLAROM 00000290
43 00000120
FOLL2 SLAROM 00000290
44 00000120
FOLL3 SLAROM 00000290
45 00000120
FOLL4 SLAROM 00000120
46 00000290
SHIEL SLAROM 00000120
47 00000290
IRDSH SLAROM 00000120
48 00000290
LSHLD SLAROM 00000120
49 00000290
RODI1 SLAROM 00000120
50 00000290
RODI2 SLAROM 00000120
51 00000290
RODI3 SLAROM 00000120
52 00000290
RODI4 SLAROM 00000120
IC01F SLAROM 00000890
023 00000300
0 0 0 7 00000910
999 00000920
00000960

```

```

HETEROGENEOUS 2 TEMPERATURE MODEL
70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
1 2
0.0 0.30705 0.60865
PIJF
37 0 5 3 0 0 0 0
0 0 0. 0. 0. 0.
0
EDIT
2 2 1 0 0
IC01FF
PREP
INNER CORE 2,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3
2 1 0 1 1 0 3 0 0 0 0 0 70
10703.15 1.3 0.0
1373.15 703.15
25 8
0.30705 0.30160
949 0.510408E-02 940 0.853884E-03 941 0.674971E-04 942 0.313286E-05
951 0.329624E-05 925 0.223986E-04 926 0.216956E-05 928 0.100072E-01
8 0.349848E-01 6 0.111864E-02
937 0.232319E-05 939 0.207934E-05 950 0.762319E-07
953 0.151805E-06 962 0.157391E-06 963 0.532589E-08 964 0.144924E-07
965 0.542844E-09 948 0.560082E-06
105 0.900654E-03
115 0.304139E-02
887 0.766799E-04 897 0.146361E-02 857 0.743652E-05
817 0.543780E-04

11 0.139463E-01 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

```

Data file [12] .OPDEND.DATA(GDI3PRT)

SLAROM input for 'perturbation'

```

PREP
INNER CORE 1,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3
2 1 0 1 1 0 3 -20 0 0 0 0 70
10703.15 1.3 0.0
1373.15 703.15
25 8
0.30705 0.30160
949 0.522699E-02 940 0.840088E-03 941 0.648861E-04 942 0.283603E-05
951 0.319672E-05 925 0.230442E-04 926 0.208237E-05 928 0.100579E-01
8 0.349848E-01 6 0.111864E-02
937 0.185217E-05 939 0.193110E-05 950 0.702519E-07
953 0.134268E-06 962 0.143732E-06 963 0.465225E-08 964 0.121468E-07
965 0.436447E-09 948 0.452596E-06
105 0.924847E-03
115 0.304139E-02
887 0.597014E-04 897 0.133347E-02 857 0.688767E-05
817 0.448337E-04

11 0.139463E-01 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

```

PATH

```

PATH
HETEROGENEOUS 2 TEMPERATURE MODEL
70 2 2 1 0 0 0
3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
1 2
0.0 0.30705 0.60865
PIJF
37 0 5 3 0 0 0 0
0 0 0. 0. 0. 0.
0
EDIT
2 2 1 0 0
IC02FF
PREP
.....
REPETITIVE DATA REMOVED
.....
OC15FF
PREP
OUTER CORE 16,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3
2 1 0 1 1 0 3 0 0 0 0 0 70
10703.15 1.3 0.0
1373.15 703.15

```

25 8
 0.30705 0.26171
 949 0.518056E-02 940 0.800749E-03 941 0.631022E-04 942 0.269689E-05
 951 0.313127E-05 925 0.189842E-04 926 0.162312E-05 928 0.811942E-02
 8 0.306073E-01 6 0.976832E-03
 937 0.128360E-05 939 0.148325E-05 950 0.672017E-07
 953 0.132617E-06 962 0.136709E-06 963 0.440933E-08 964 0.119792E-07
 965 0.440346E-09 948 0.354594E-06
 105 0.821440E-03
 115 0.265585E-02
 887 0.407280E-04 897 0.118156E-02 857 0.516368E-05
 817 0.378454E-04

11 0.146730E-01 26 0.166387E-01 24 0.463825E-02 28 0.326322E-02
 42 0.369672E-03 25 0.387342E-03 40 0.180912E-02 1 0.307549E-02

PATH

HETEROGENEOUS 2 TEMPERATURE MODEL

70 2 2 1 0 0 0
 3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0

1 2
 0.0 0.30705 0.56876

PIJF

37 0 5 3 0 0 0 0
 0 0 0. 0. 0. 0.

EDIT

2 2 1 0 0

OC16FF

PREP

PLENUM ,LO Q ,FU,165%FUEL, UPPER INNER

1 1 0 0 0 0 3 0 0 0 0 0 0 70
 703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
 42 2.61950E-04 25 2.74470E-04

UIPLANN

PREP

PLUGGED PIN, LO Q ,165%FUEL UPPER INNER

1 1 0 0 0 0 3 0 0 0 0 0 0 70
 703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 2.96650E-02 24 8.26930E-03 28 5.81800E-03
 42 6.59100E-04 25 6.90610E-04

UIPLGG

PREP

ROD FOLLOWER 1

1 1 0 0 0 0 3 0 0 0 0 0 0 70
 703.15 0.0 0.0

6

1.0

11 1.99080E-02 26 5.24170E-03 24 1.46120E-03 28 1.02800E-03
 42 1.16460E-04 25 1.22020E-04

FOLLW

PREP

UPPER AXIAL & OUTER RADIAL SHIELD

1 1 0 0 0 0 3 0 0 0 0 0 0 70

703.15 0.0 0.0

6

1.0

11 4.38990E-03 26 4.50900E-02 24 1.25690E-02 28 8.84310E-03
 42 1.00180E-03 25 1.04970E-03

SHIELD

PREP

LOWER SHIELD,

1 1 0 0 0 0 3 0 0 0 0 0 70

703.15 0.0 0.0

9

1.0

11 6.36540E-03 26 1.77540E-02 24 4.94900E-03 28 3.48200E-03
 42 3.94460E-04 25 4.13320E-04 6 9.24600E-03 105 7.89610E-03
 115 2.90880E-02

LSHDD

PREP

ABSORBER ROD 1, HETEROGENEOUS CELL MODEL WITHIN 9.7 S/A FUEL

6 1 0 1 1 0 3 0 0 0 0 0 70

703.15 1.3 0.0

3 1 5 1 5 30

4.7901 0.1137 0.9751 2.1419 0.3937 18.8519

6 2.47080E-02 105 9.15890E-02 115 7.24350E-03 /

11 2.19495E-09 /

26 5.63630E-02 24 1.57110E-02 28 1.10540E-02

42 1.25230E-03 25 1.31210E-03 /

11 2.19495E-02 /

26 5.63630E-02 24 1.57110E-02 28 1.10540E-02

42 1.25230E-03 25 1.31210E-03 /

949 6.49311E-04 940 8.17696E-04 941 1.68546E-04 942 3.13512E-04

951 1.94612E-05 925 7.78987E-06 926 3.61466E-07 928 2.99340E-03

8 1.03332E-02 11 1.18383E-02 26 1.07946E-02 24 3.00913E-03

28 2.11706E-03 42 2.39830E-04 25 2.51294E-04

40 6.79872E-04

937 3.26805E-07 939 6.64538E-07 950 4.94608E-07 953 1.17133E-05

962 1.13309E-06 963 4.51328E-08 964 1.29052E-06 965 4.89657E-08

948 6.45035E-05

1 1.15578E-03

887 1.02518E-05 897 8.98626E-05 857 1.21029E-06

817 6.73348E-05 /

PATH

TITLE CARD HETEROGENEOUS ROD

70 6 2 1 1 0 0

3 6 6 0 6 0 0 0 0 0 8 10 0 0 15 0

1 2 3 4 5 6

0.0 4.7901 4.9038 5.7652 7.9071 8.3008 27.1527

PIJF

37 0 5 3 0 1 0 0

0 0 0. 0. 0. 0.

0

EDIT

2 2 1 1 0

1 5 1 5

ROD11A

PREP

INNER CORE 1,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3 NA VOID

2 1 0 1 1 0 3 -20 0 0 0 0 70

10703.15 1.3 0.0

1373.15 703.15
 25 8
 0.30705 0.30160
 949 0.522699E-02 940 0.840088E-03 941 0.648861E-04 942 0.283603E-05
 951 0.319672E-05 925 0.230442E-04 926 0.208237E-05 928 0.100579E-01
 8 0.349848E-01 6 0.111864E-02
 937 0.185217E-05 939 0.193110E-05 950 0.702519E-07
 953 0.134268E-06 962 0.143732E-06 963 0.465225E-08 964 0.121468E-07
 965 0.436447E-09 948 0.452596E-06
 105 0.924847E-03
 115 0.304139E-02
 887 0.597014E-04 897 0.133347E-02 857 0.688767E-05
 817 0.448337E-04

 11 0.000000E-00 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
 42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

PATH
 HETEROGENEOUS 2 TEMPERATURE MODEL
 70 2 2 1 0 0 0
 3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
 1 2
 0.0 0.30705 0.60865
 PIJF
 37 0 5 3 0 0 0 0
 0 0 0. 0. 0. 0.
 0
 EDIT
 2 2 1 0 0
 IC01VV
 PREP

REPETITIVE DATA REMOVED

OC15VV
 PREP
 OUTER CORE 16,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3
 2 1 0 1 1 0 3 0 0 0 0 0 70
 10703.15 1.3 0.0
 1373.15 703.15
 25 8
 0.30705 0.26171
 949 0.518056E-02 940 0.800749E-03 941 0.631022E-04 942 0.269689E-05
 951 0.313127E-05 925 0.189842E-04 926 0.162312E-05 928 0.811942E-02
 8 0.306073E-01 6 0.976832E-03
 937 0.128360E-05 939 0.148325E-05 950 0.672017E-07
 953 0.132617E-06 962 0.136709E-06 963 0.440933E-08 964 0.119792E-07
 965 0.440346E-09 948 0.354594E-06
 105 0.821440E-03
 115 0.265585E-02
 887 0.407280E-04 897 0.118156E-02 857 0.516368E-05
 817 0.378454E-04

 11 0.000000E-00 26 0.166387E-01 24 0.463825E-02 28 0.326322E-02
 42 0.369672E-03 25 0.387342E-03 40 0.180912E-02 1 0.307549E-02

PATH
 HETEROGENEOUS 2 TEMPERATURE MODEL
 70 2 2 1 0 0 0
 3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
 1 2
 0.0 0.30705 0.56876
 PIJF
 37 0 5 3 0 0 0 0
 0 0 0. 0. 0. 0.
 0
 EDIT
 2 2 1 0 0
 OC16VV
 PREP
 INNER CORE 1,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3 DOPPLER
 2 1 0 1 1 0 3 -20 0 0 0 0 70
 10703.15 1.3 0.0
 1873.15 703.15
 25 8
 0.30705 0.30160
 949 0.522699E-02 940 0.840088E-03 941 0.648861E-04 942 0.283603E-05
 951 0.319672E-05 925 0.230442E-04 926 0.208237E-05 928 0.100579E-01
 8 0.349848E-01 6 0.111864E-02
 937 0.185217E-05 939 0.193110E-05 950 0.702519E-07
 953 0.134268E-06 962 0.143732E-06 963 0.465225E-08 964 0.121468E-07
 965 0.436447E-09 948 0.452596E-06
 105 0.924847E-03
 115 0.304139E-02
 887 0.597014E-04 897 0.133347E-02 857 0.688767E-05
 817 0.448337E-04

 11 0.139463E-01 26 0.158146E-01 24 0.440852E-02 28 0.310159E-02
 42 0.351362E-03 25 0.368157E-03 40 0.171951E-02 1 0.292317E-02

PATH
 HETEROGENEOUS 2 TEMPERATURE MODEL
 70 2 2 1 0 0 0
 3 2 2 0 2 0 0 0 4 0 6 180 0 0 360 0
 1 2
 0.0 0.30705 0.60865
 PIJF
 37 0 5 3 0 0 0 0
 0 0 0. 0. 0. 0.
 0
 EDIT
 2 2 1 0 0
 IC01DD
 PREP

REPETITIVE DATA REMOVED

OC15DD
 PREP
 OUTER CORE 16,C7,HPU40/45%,4*9.75M,165%,24/8ZRH/B4C,PERT,GDI3
 2 1 0 1 1 0 3 0 0 0 0 70
 10703.15 1.3 0.0


```

USERPDS OC12FF CITA 00000080
INNER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN OC13FF SLAR 0 1 00000060
PDSOUT OC13FF SLAR 00000070
USERPDS OC13FF CITA 00000080
INNER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN OC14FF SLAR 0 1 00000060
PDSOUT OC14FF SLAR 00000070
USERPDS OC14FF CITA 00000080
INNER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN OC15FF SLAR 0 1 00000060
PDSOUT OC15FF SLAR 00000070
USERPDS OC15FF CITA 00000080
INNER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN OC16FF SLAR 0 1 00000060
PDSOUT OC16FF SLAR 00000070
USERPDS OC16FF CITA 00000080
INNER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN UIPLNN SLAR 0 1 00000060
PDSOUT UIPLNN SLAR 00000070
USERPDS UIPLNN CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN UIPLGG SLAR 0 1 00000060
PDSOUT UIPLGG SLAR 00000070
USERPDS UIPLGG CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN FOLL1W SLAR 0 1 00000060
PDSOUT FOLL1W SLAR 00000070
USERPDS FOLL1W CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN SHIELD SLAR 0 1 00000060
PDSOUT SHIELD SLAR 00000070
USERPDS SHIELD CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN LSHLDD SLAR 0 1 00000060
PDSOUT LSHLDD SLAR 00000070
USERPDS LSHLDD CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020
PDSIN RODILA SLAR 0 1 00000060
PDSOUT RODILA SLAR 00000070
USERPDS RODILA CITA 00000080
OUTER-CORE 182*4DAY,HI Q PU , 40% T.D. HAA ITER. 1 00000020

```

```

1 0 0 1 1 0 0 1 1 2 0 0 0 1 0 0 0 0 0 0 0 0 0 0 0 0
003
0 0 0 0 7 0 0 0 0 0 1 0 0 0 0 0 0 1 0 0 0 0 0
5.0000E-6 1.0000E-6 1.5600E+3
004
10 46.21720 1 4.27490 7 38.52480 1 2.29260 1 2.23650 7 38.48690
1 4.61620 6 28.74400 1 1.24510 1 2.46280 5 21.29210 2 6.00000
1 3.00000 4 21.10000 4 21.10000
10 50.00000 2 7.00000 1 3.00000 2 5.00000 1 3.00000 2 6.00000
3 9.00000 4 12.00000 4 12.00000 3 9.00000 2 6.00000 1 3.00000
1 2.00000 14 70.00000 6 33.00000
005
20 22 20 22 22 20 22 20 22 22 20 20 20 20 20
19 22 19 22 22 19 22 19 22 22 19 19 19 19 20 20
18 22 18 22 22 18 22 18 22 22 18 18 18 18 20 20
17 22 17 22 22 17 22 17 22 22 17 17 17 17 20 20
1 19 1 19 19 1 19 9 19 19 9 9 9 20 20
2 19 2 19 19 2 19 10 19 19 10 10 10 20 20
3 19 3 19 19 3 19 11 19 19 11 11 11 20 20
4 19 4 19 19 4 19 12 19 19 12 12 12 20 20
5 19 5 19 19 5 19 13 19 19 13 13 13 20 20
6 19 6 19 19 6 19 14 19 19 14 14 14 20 20
7 19 7 19 19 7 19 15 19 19 15 15 15 20 20
8 19 8 19 19 8 19 16 19 19 16 16 16 20 20
18 19 18 19 19 18 19 18 19 19 18 18 18 20 20
17 19 17 19 19 17 19 17 19 19 17 17 17 20 20
21 21 21 21 21 21 21 21 21 21 21 21 21 20 20
008
-70 70 0 22 2 1 1 3
IC01VV SLAROM 00000290
2 00000120
IC02VV SLAROM 00000290
3 00000120
IC03VV SLAROM 00000290
4 00000120
IC04VV SLAROM 00000290
5 00000120
IC05VV SLAROM 00000290
6 00000120
IC06VV SLAROM 00000290
7 00000120
IC07VV SLAROM 00000290
8 00000120
IC08VV SLAROM 00000290
9 00000290
OC09VV SLAROM 00000120
10 00000290
OC10VV SLAROM 00000120
11 00000290
OC11VV SLAROM 00000120
12 00000290
OC12VV SLAROM 00000120
13 00000290
OC13VV SLAROM 00000120
14 00000290

```

Data file [16] .OPCITD.DATA(PIJ3PV70)

CITATION input, void, 70 groups, for 'perturbation'

```

CITATION 00000010
IC01VV SLAROM 00000120
600MW / 60 CM PU BURNING CORE , PU VECTOR STUDY , LOW QUALITY PU
PERTURBATION 70 GROUP VOID , CORE 7 , 165% PIN SIZE, RODS 0% IN
001
0 0 0 0 0 0 1 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 1 0 0 0 0 0

```



```

IC02FF SLAROM 00000120
IC03FF SLAROM 00000120
IC04FF SLAROM 00000120
IC05FF SLAROM 00000120
IC06FF SLAROM 00000120
IC07FF SLAROM 00000120
IC08FF SLAROM 00000120
OC09FF SLAROM 00000120
OC10FF SLAROM 00000120
OC11FF SLAROM 00000120
OC12FF SLAROM 00000120
OC13FF SLAROM 00000120
OC14FF SLAROM 00000120
OC15FF SLAROM 00000120
OC16FF SLAROM 00000120
UIPLNN SLAROM 00000120
UIPLGG SLAROM 00000120
FOLLW SLAROM 00000120
SHIELD SLAROM 00000120
LSHLDD SLAROM 00000120
RODILA SLAROM 00000120
004 00000910
2 00000920
006 00000960
0.00000E+00 0.25000E+00 0.50000E+00 0.75000E+00 1.00000E+00 1.25000E+00 000000920
1.50000E+00 1.75000E+00 2.00000E+00 2.25000E+00 2.50000E+00 2.75000E+00 000000920
3.00000E+00 3.25000E+00 3.50000E+00 3.75000E+00 4.00000E+00 4.25000E+00 000000920
4.50000E+00 4.75000E+00 5.00000E+00 5.25000E+00 5.50000E+00 5.75000E+00 000000920
6.00000E+00 6.25000E+00 6.50000E+00 6.75000E+00 7.00000E+00 7.25000E+00 000000920
7.50000E+00 7.75000E+00 8.00000E+00 8.25000E+00 8.50000E+00 8.75000E+00 000000920
9.00000E+00 9.25000E+00 9.50000E+00 9.75000E+00 1.00000E+01 1.02500E+01 000000920
1.05000E+01 1.07500E+01 1.10000E+01 1.12500E+01 1.15000E+01 1.17500E+01 0100000920
1.20000E+01 1.22500E+01 1.25000E+01 1.27500E+01 1.30000E+01 1.32500E+01 0100000920
1.35000E+01 1.37500E+01 1.40000E+01 1.42500E+01 1.45000E+01 1.47500E+01 0100000920
1.50000E+01 1.52500E+01 1.55000E+01 1.57500E+01 1.60000E+01 1.62500E+01 0100000920
1.65000E+01 1.67500E+01 1.70000E+01 2.76500E+01 00000920
999 00000960
00000960

```

Data file [20] .OPEDID.DATA(PIJ3PDNX)

PERKY input, delayed neutron fraction, for 'perturbation'

```

PERKY 00000010
IC01FF SLAROM 00000120
18 GROUP DELAYED NEUTRON FRACTION
001 1 18 22 1 0 0 0 1
002 1 0 0
16 16 0 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16
1 8
IC01FF SLAROM 00000270

```

```

949 940 941 942 925 928 948 951
2 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
3 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
4 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
5 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
6 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
7 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
8 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
9 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
10 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
11 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
12 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
13 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
14 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
15 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
16 8
IC01FF SLAROM 00000270
949 940 941 942 925 928 948 951
1 1 2 2 3 3 4 4 5 5 6 6 7 7 8 8 9 9 10 10 11 11 12 12
13 13 14 14 15 15 16 16 17 18 19 20 21 22
003
1 0 0
22 22 0 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20
21 22
IC01FF SLAROM 00000270
18 18 0 2 0 0 1 1 1
IC01FF SLAROM 00000120
IC02FF SLAROM 00000120
IC03FF SLAROM 00000120
IC04FF SLAROM 00000120
IC05FF SLAROM 00000120

```



```

UIPLGG SLAROM 00000120
FOLL1W SLAROM 00000120
SHIELD SLAROM 00000120
LSHLDD SLAROM 00000120
ROD11A SLAROM 00000120
IC01VV SLAROM 00000120
IC02VV SLAROM 00000120
IC03VV SLAROM 00000120
IC04VV SLAROM 00000120
IC05VV SLAROM 00000120
IC06VV SLAROM 00000120
IC07VV SLAROM 00000120
IC08VV SLAROM 00000120
OC09VV SLAROM 00000120
OC10VV SLAROM 00000120
OC11VV SLAROM 00000120
OC12VV SLAROM 00000120
OC13VV SLAROM 00000120
OC14VV SLAROM 00000120
OC15VV SLAROM 00000120
OC16VV SLAROM 00000120
004 00000910
5 00000920
008 00000960
1 0 1 16 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 00000920
IC01VV SLAROM 00000960
1 23 2 24 3 25 4 26 5 27 6 28 7 29 8 30 9 31 10 32 11 33 12 34 00000960
13 35 14 36 15 37 16 38 00000960
999 00000960

```

Data file [22] .OPEDID.DATA(PIJ3PFD0)

PERKY input, Doppler (by zone), for 'perturbation'

```

PERKY 00000010
IC01FF SLAROM 00000120
18 GROUP DOPPLER PERTURBATION, FIRST ORDER, BY REGION
001 1 18 38 1 0 0 0 1
003 1 0 0
38 38 0 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20
21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38
IC01FF SLAROM 00000270
18 18 0 2 0 0 1 1 1 00000280
IC01FF SLAROM 00000120
IC02FF SLAROM 00000120
IC03FF SLAROM 00000120
IC04FF SLAROM 00000120
IC05FF SLAROM 00000120
IC06FF SLAROM 00000120
IC07FF SLAROM 00000120

```

```

IC08FF SLAROM 00000120
OC09FF SLAROM 00000120
OC10FF SLAROM 00000120
OC11FF SLAROM 00000120
OC12FF SLAROM 00000120
OC13FF SLAROM 00000120
OC14FF SLAROM 00000120
OC15FF SLAROM 00000120
OC16FF SLAROM 00000120
UIPLNN SLAROM 00000120
UIPLGG SLAROM 00000120
FOLL1W SLAROM 00000120
SHIELD SLAROM 00000120
LSHLDD SLAROM 00000120
ROD11A SLAROM 00000120
IC01DD SLAROM 00000120
IC02DD SLAROM 00000120
IC03DD SLAROM 00000120
IC04DD SLAROM 00000120
IC05DD SLAROM 00000120
IC06DD SLAROM 00000120
IC07DD SLAROM 00000120
IC08DD SLAROM 00000120
OC09DD SLAROM 00000120
OC10DD SLAROM 00000120
OC11DD SLAROM 00000120
OC12DD SLAROM 00000120
OC13DD SLAROM 00000120
OC14DD SLAROM 00000120
OC15DD SLAROM 00000120
OC16DD SLAROM 00000120
004 00000910
4 00000920
008 00000960
1 0 1 16 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 00000920
IC01DD SLAROM 00000960
1 23 2 24 3 25 4 26 5 27 6 28 7 29 8 30 9 31 10 32 11 33 12 34 00000960
13 35 14 36 15 37 16 38 00000960
999 00000960

```

Data file [23] .OPEDID.DATA(GDI3PFD1)

PERKY input, Doppler (by nuclide), for 'perturbation'

```

PERKY 00000010
IC01FF SLAROM 00000120
18 GROUP DOPPLER PERTUREATION, 1'ST ORDER, BY ISOTOPE, 1 TO 10 (OF 20)
001 1 18 38 1 0 0 0 1
002 1 0 0
38 38 0 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20

```



```

40 1
32 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
33 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
34 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
35 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
36 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
37 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
38 20 00000280
IC01DD SLAROM 00000120
949 940 941 942 951 925 926 928 8 11 26 24 28 42 6 953 115 10500000120
40 1
1 1 2 2 3 3 4 4 5 5 6 6 7 7 8 8 9 9 10 10 11 11 12 12
13 13 14 14 15 15 16 16 17 17 18 18 19 19 20 20 21 21 22 22 23 23 24 24
25 25 26 26 27 27 28 28 29 29 30 30 31 31 32 32 33 33 34 34 35 35 36 36
37 37 38 38
003
1 0 0
38 38 0 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20
21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38
IC01FF SLAROM 00000270
18 18 0 2 0 0 1 1 1 00000280
IC01FF SLAROM 00000120
IC02FF SLAROM 00000120
IC03FF SLAROM 00000120
IC04FF SLAROM 00000120
IC05FF SLAROM 00000120
IC06FF SLAROM 00000120
IC07FF SLAROM 00000120
IC08FF SLAROM 00000120
OC09FF SLAROM 00000120
OC10FF SLAROM 00000120
OC11FF SLAROM 00000120
OC12FF SLAROM 00000120
OC13FF SLAROM 00000120
OC14FF SLAROM 00000120
OC15FF SLAROM 00000120
OC16FF SLAROM 00000120
UIPLNN SLAROM 00000120
UIPLGG SLAROM 00000120
FOLL1W SLAROM 00000120
SHIELD SLAROM 00000120

```

```

LSHLDD SLAROM 00000120
ROD1IA SLAROM 00000120
IC01DD SLAROM 00000120
IC02DD SLAROM 00000120
IC03DD SLAROM 00000120
IC04DD SLAROM 00000120
IC05DD SLAROM 00000120
IC06DD SLAROM 00000120
IC07DD SLAROM 00000120
IC08DD SLAROM 00000120
OC09DD SLAROM 00000120
OC10DD SLAROM 00000120
OC11DD SLAROM 00000120
OC12DD SLAROM 00000120
OC13DD SLAROM 00000120
OC14DD SLAROM 00000120
OC15DD SLAROM 00000120
OC16DD SLAROM 00000120
004 00000910
4 00000920
008 00000960
0 1 1 16 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 00000920
IC01DD SLAROM 00000120
10 949 940 941 942 951 925 926 928 8 11 00000960
949 940 941 942 951 925 926 928 105 115 00000960
1 23 10 00000960
949 0.0 0.0 00000960
940 0.0 0.0 00000960
941 0.0 0.0 00000960
942 0.0 0.0 00000960
951 0.0 0.0 00000960
925 0.0 0.0 00000960
926 0.0 0.0 00000960
928 0.0 0.0 00000960
8 0.0 0.0 00000960
11 0.0 0.0 00000960
2 24 10 00000960
949 0.0 0.0 00000960
940 0.0 0.0 00000960
941 0.0 0.0 00000960
942 0.0 0.0 00000960
951 0.0 0.0 00000960
925 0.0 0.0 00000960
926 0.0 0.0 00000960
928 0.0 0.0 00000960
8 0.0 0.0 00000960
11 0.0 0.0 00000960
3 25 10 00000960
949 0.0 0.0 00000960
940 0.0 0.0 00000960
941 0.0 0.0 00000960
942 0.0 0.0 00000960
951 0.0 0.0 00000960
925 0.0 0.0 00000960
926 0.0 0.0 00000960
928 0.0 0.0 00000960
8 0.0 0.0 00000960
11 0.0 0.0 00000960
4 26 10 00000960

```



```

      8 0.0      0.0      00000960
     11 0.0      0.0      00000960
    15 37 10      0.0      00000960
     949 0.0      0.0      00000960
     940 0.0      0.0      00000960
     941 0.0      0.0      00000960
     942 0.0      0.0      00000960
     951 0.0      0.0      00000960
     925 0.0      0.0      00000960
     926 0.0      0.0      00000960
     928 0.0      0.0      00000960
      8 0.0      0.0      00000960
     11 0.0      0.0      00000960
    16 38 10      0.0      00000960
     949 0.0      0.0      00000960
     940 0.0      0.0      00000960
     941 0.0      0.0      00000960
     942 0.0      0.0      00000960
     951 0.0      0.0      00000960
     925 0.0      0.0      00000960
     926 0.0      0.0      00000960
     928 0.0      0.0      00000960
      8 0.0      0.0      00000960
     11 0.0      0.0      00000960
999

```

Data file [24] .OPMOSD.DATA(GDI1DY)

MOSES input, steady state, for 'rod worth'

GDI1S-HQPU40.25/45.03%,165%F,4*9.75M.,24/8PZRH/B4C(30%B10),45%HOLE,OPT.
 SOC, PCR 50%, BCR 100% BCR Y(404) ROD STUCK B OK 7 GROUPS
 && CORE 7, DETAILED, IMPROVED AXIAL REFLECTORS RODS HOMO

```

&&
&&
CONTROL
  3*1      && DIFFUSION CAL. ON
  3*0      && BURNUP CAL. ON
EDIT
  0 4*1 2 19*0 1 4*0      && EDIT PROGRAM SETUP
  0 1 3*0 1 3*0 1 20*0      && EDIT FUEL MANAGEMENT DATA
  0 0 2 2 1 0 1 0 1 1 0 1 0 1 3 0 0 1 1 0 0 1 3*0 1 0 1 0 0      && DIF CAL
  3*0 1 4*0 1 0 2 5*0 9*0 1 4*0      && EDIT BURNUP CAL. DATA

```

```

EDITD
&& $IC1 $IC2 $OC1
&& DETAILED EDIT FOR DIFF. CAL.
  50*0.0
&& DETAILED EDIT FOR BURNUP CAL.
  50*0.0
&& DETAILED EDIT FOR N.D. CAL.
  50*0.0
EOB

```

```

&&
&&
&&
LOADING 1
&&
GEOM
&& NSYM NLAY NPLN NAZN NGRP PITCH
      6 18 56 15 7 15.81
&& AXIAL MESH BY REGION
      10 2 1 2 1 2 3 4 4 3 2 1 1 14 6
&& REGION WIDTH (CM)
      50.0 7.0 3.0 5.0 3.0 6.0 9.0 12.0 12.0 9.0 6.0 3.0 2.0 70.0 33.0
&& === NMES ( NMES=1:CM, NMES=2:MOD.CM)
&& NMES NCAL NITR NIIT NITM NBATCH OUTCR RLOSS DELAY NSMALL
CPARA1 2 0 200 3 1600 0 0 0 0 0 3*0 182.5 1
CPARA2 5.E-6 1.E-4 0.5 1.5 1 0.0 4*0 0.0 0.0
EOB
&&
&&

```

```

&& ASSEMBLY AXIAL INFORMATION
AXIAL
&& NREGZ
15 28 27 26 25 1 2 3 4 5 6 7 8 26 25 29 && INNER FUEL
    50.0 7.0 3.0 5.0 3.0 6.0 9.0 12.0 12.0 9.0 6.0 3.0 2.0 70.0 33.0
15 28 27 26 25 9 10 11 12 13 14 15 16 26 25 29 && OUTER FUEL1
    50.0 7.0 3.0 5.0 3.0 6.0 9.0 12.0 12.0 9.0 6.0 3.0 2.0 70.0 33.0
15 28 27 26 25 17 18 19 20 21 22 23 24 26 25 29 && OUTER FUEL2
    50.0 7.0 3.0 5.0 3.0 6.0 9.0 12.0 12.0 9.0 6.0 3.0 2.0 70.0 33.0
1 28 && RADIAL SHIELD
    230.0
3 30 27 29 && ROD 0%
    65.0 132.0 33.0
4 27 30 27 29 && ROD 50%
    50.0 45.0 102.0 33.0
4 27 30 27 29 && ROD 100%
    65.0 60.0 72.0 33.0

```

```

EOI
&& ASSEMBLY GROUP ASSIGNMENT
ASYGROUP
  1 1 1 1 1 IC1 INNER CORE
  1 2 1 2 2 OC1 OUTER CORE1
  1 3 1 2 2 OC2 OUTER CORE2
  1 4 1 4 3 SHD RADIAL SHIELD
  2 5 1 4 -4 R0% ROD 0%
  2 6 1 4 -5 R50 ROD 50%
  2 7 1 4 -6 RIN ROD 100%
EOI
&& MATERIAL ASSIGNMENT
REGION
  1 1 1 1 1 $IC1 INNER CORE1
  1 1 1 1 1 $IC2 INNER CORE2
  1 1 1 1 1 $IC3 INNER CORE3
  1 1 1 1 1 $IC4 INNER CORE4
  1 1 1 1 1 $IC5 INNER CORE5
  1 1 1 1 1 $IC6 INNER CORE6
  1 1 1 1 1 $IC7 INNER CORE7
  1 1 1 1 1 $IC8 INNER CORE8
  2 2 1 1 1 $O11 OUTER CORE1/1

```

2	2	1	1	1	\$O12	OUTER CORE1/2
2	2	1	1	1	\$O13	OUTER CORE1/3
2	2	1	1	1	\$O14	OUTER CORE1/4
2	2	1	1	1	\$O15	OUTER CORE1/5
2	2	1	1	1	\$O16	OUTER CORE1/6
2	2	1	1	1	\$O17	OUTER CORE1/7
2	2	1	1	1	\$O18	OUTER CORE1/8
2	3	1	1	1	\$O21	OUTER CORE2/1
2	3	1	1	1	\$O22	OUTER CORE2/2
2	3	1	1	1	\$O23	OUTER CORE2/3
2	3	1	1	1	\$O24	OUTER CORE2/4
2	3	1	1	1	\$O25	OUTER CORE2/5
2	3	1	1	1	\$O26	OUTER CORE2/6
2	3	1	1	1	\$O27	OUTER CORE2/7
2	3	1	1	1	\$O28	OUTER CORE2/8
3	4	1	3	1	\$PLN	PLENUM AXIAL
4	5	1	3	1	\$PLG	PLUG AXIAL
5	6	1	3	1	\$FOL	GAP FOLLOWER
6	7	1	5	1	\$URS	UP/RAD SHLD
7	8	1	3	1	\$LSH	LOWER SHIELD
9	9	1	6	1	\$ABS	ROD ABSORBER

EOI
 &&
 && CROSS SECTION AND FISSION SPECTRUM
 &&

MICROXS
 1 9 33 2 1 2
 1.4782E-1 4.5751E-1 3.8563E-1 8.7542E-3 2.807E-4 0.0 0.0

&&
 && NUCLIDE INFORMATION

CODE	NAME	IFIS	KIND	BCAL	KLIB	ATW	EFIS	ECAP	DECAY
1	PU239	2	1	0	1	239.05	3.34E-11	0	0
2	PU240	2	2	0	1	240.05	3.36E-11	0	0
3	PU241	2	1	0	1	241.06	3.37E-11	0	1.526E-9
4	PU242	2	3	0	1	242.06	3.38E-11	0	0
5	AM241	2	3	0	1	241.06	3.36E-11	0	0
6	U235	1	3	0	1	235.044	3.23E-11	0	0
7	U236	1	3	0	1	236.05	3.24E-11	0	0
8	U238	1	2	0	1	238.051	3.31E-11	0	0
9	O	0	5	0	1	15.999	0	0	0
10	NA	0	6	0	1	22.990	0	0	0
11	FE	0	6	0	1	55.847	0	0	0
12	CR	0	6	0	1	51.996	0	0	0
13	NI	0	6	0	1	58.70	0	0	0
14	MO	0	6	0	1	95.94	0	0	0
15	MN	0	6	0	1	54.938	0	0	0
18	C	0	7	0	1	12.011	0	0	0
20	ZR	0	6	0	1	91.224	0	0	0
21	NP237	0	6	0	1	237.048	3.30E-11	0	0
22	NP239	0	6	0	1	239.053	3.31E-11	0	0
23	AM242M	0	6	0	1	242.060	3.45E-11	0	0
24	AM243	0	6	0	1	243.061	3.46E-11	0	0
25	CM242	0	6	0	1	242.059	3.52E-11	0	0
26	CM243	0	6	0	1	243.060	3.52E-11	0	0
27	CM244	0	6	0	1	244.063	3.53E-11	0	0
28	CM245	0	6	0	1	245.066	3.53E-11	0	0
29	PU238	2	3	0	1	238.05	3.33E-11	0	0

67	E-10	0	7	0	1	10.0123	0	0	0
68	B-11	0	7	0	1	11.0093	0	0	0
70	H	0	7	0	1	1.0078	0	0	0
78	U238-FP	-1	4	0	1	238.051	0	0	0
79	PU239-FP	-1	4	0	1	239.055	0	0	0
80	U235-FP	-1	4	0	1	235.044	0	0	0
81	PU241-FP	-1	4	0	1	241.057	0	0	0

REACTI

6	7	2	1.0
8	1	2	1.0
1	2	2	1.0
2	3	2	1.0
3	4	2	1.0
29	1	2	1.0
3	5	1	1.0

YIELD

78	8	1.0
79	1	1.0
79	2	1.0
79	29	1.0
80	6	1.0
80	7	1.0
81	3	1.0
81	4	1.0
81	5	1.0

&&
 && ATOM NUMBER DENSITY
 &&

ATDEN

&& FROM .OPFORT(GDI1DRZ)

&& ZONE 1

1.46918E-03	1.74999E-04	9.65638E-06	3.30258E-07
3.79707E-07	6.73607E-06	3.35554E-07	2.60913E-03
8.90358E-03	1.03968E-02	1.17896E-02	3.28651E-03
2.31220E-03	2.61937E-04	2.74457E-04	
	2.84691E-04		1.28188E-03
2.91854E-07	3.71285E-07	6.62542E-09	1.22009E-08
1.51758E-08	3.82196E-10	8.50733E-10	2.37739E-11
4.90057E-08			
2.83107E-04	7.74029E-04		
2.17919E-03			
9.15262E-06	2.09324E-04	1.09239E-06	
5.96565E-06			

&& ZONE 2

1.44892E-03	1.77607E-04	1.00720E-05	3.65641E-07
3.93013E-07	6.62858E-06	3.51142E-07	2.60133E-03
8.90358E-03	1.03968E-02	1.17896E-02	3.28651E-03
2.31220E-03	2.61937E-04	2.74457E-04	
	2.84691E-04		1.28188E-03
3.68157E-07	4.00145E-07	7.22962E-09	1.38310E-08
1.66777E-08	4.39829E-10	1.01782E-09	2.96814E-11
6.21125E-08			
2.78976E-04	7.74029E-04		
2.17919E-03			
1.17643E-05	2.30406E-04	1.18318E-06	
7.26002E-06			

&& ZONE 3

REPETITIVE DATA REMOVED

```

&& ZONE 24
1.72011E-03 1.77176E-04 7.91079E-06 1.98700E-07
3.22824E-07 6.58690E-06 2.35864E-07 2.42671E-03
8.92029E-03 1.03968E-02 1.17896E-02 3.28651E-03
2.31221E-03 2.61937E-04 2.74458E-04
      2.84691E-04      1.28188E-03
1.49815E-07 2.43232E-07 4.18611E-09 5.90895E-09
9.29809E-09 1.76758E-10 3.03483E-10 6.46939E-12
2.18182E-08
3.02515E-04 7.74028E-04
2.17919E-03
4.48776E-06 1.57107E-04 7.08526E-07
3.33516E-06
&& ZONE 25
0 0 0 0 0 0 0 0
1.03970E-02 1.17900E-02 3.28650E-03 2.31220E-03
2.61950E-04 2.74470E-04
0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
&& ZONE 26
0 0 0 0 0 0 0 0
1.03970E-02 2.96650E-02 8.26929E-03 5.81799E-03
6.59100E-04 6.90609E-04
0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
&& ZONE 27
0 0 0 0 0 0 0 0
1.99080E-02 5.24169E-03 1.46120E-03 1.02800E-03
1.16460E-04 1.22020E-04
0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
&& ZONE 28
0 0 0 0 0 0 0 0
4.38990E-03 4.50900E-02 1.25690E-02 8.84310E-03
1.00180E-03 1.04970E-03
0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
&& ZONE 29
0 0 0 0 0 0 0 0
6.36540E-03 1.77540E-02 4.94900E-03 3.48200E-03
3.94460E-04 4.13320E-04 9.24600E-03 0 0 0 0 0 0 0 0 0 0 7.89610E-03
2.90880E-02
0 0 0 0 0
&& ZONE 30
0 0 0 0 0 0 0 0
9.32849E-03 1.27380E-02 3.55070E-03 2.49820E-03
2.83010E-04 2.96540E-04 8.22779E-03 0 0 0 0 0 0 0 0 0 0 3.04990E-02
2.41210E-03
0 0 0 0 0
EOI
&& === CAUTION === REFLECTIVE B.C. IS NOT PERMITTED
&& 0.0 MEANS VACUUM
BOUNDARY
7*0.0
7*0.0
7*0.0
REFUEL
1 0 365.0

```

```

0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
0 0 0.0
REFMAP
1STSCORE
      1 271 1
      272 547 2
      548 721 4
      553 557 3
      567 571 3
      581 585 3
      595 599 3
      609 613 3
      623 627 3
      547 547 3
      470 471 3
      482 484 3
      495 497 3
      508 510 3
      521 523 3
      534 536 3
      722 817 4
      825 829 4
      842 846 4
      859 863 4
      876 880 4
      893 897 4
      910 914 4
      92 92 7
      98 98 7
      104 104 7
      110 110 7
      116 116 7
      122 122 7
      404 404 5
      416 416 7
      428 428 7
      440 440 7
      452 452 7
      464 464 7
      20 20 6
      23 23 6
      26 26 6
      29 29 6
      32 32 6
      35 35 6
      95 95 6
      101 101 6
      107 107 6
      113 113 6
      119 119 6
      125 125 6
      218 218 6

```

221 221 6
 224 224 6
 227 227 6
 230 230 6
 233 233 6
 236 236 6
 239 239 6
 242 242 6
 245 245 6
 248 248 6
 251 251 6
 254 254 6
 257 257 6
 260 260 6
 263 263 6
 266 266 6
 269 269 6
 401 401 6
 407 407 6
 413 413 6
 419 419 6
 425 425 6
 431 431 6
 437 437 6
 443 443 6
 449 449 6
 455 455 6
 461 461 6
 467 467 6
 0 0 0

EOI
 PLANT
 1560.0 380
 CRPOS
 && IC1 OCl RS1 MCR BCR CRP
 0.0 0.0 0.0 0.0 0.0 0.0 34*0.0
 EOB

Data file [25] .OPDEND.DATA(GDI1RZEC)

SLAROM input for 'shutdown margin'

PREP
 INNER CORE, GDI1 (OPTIMIZED) EXPANSION COEFFICIENT
 1 1 0 0 0 0 0 3 -20 0 0 0 0 70
 1373.15 0.0 0.0
 85
 1.0
 949 1.40054E-03 940 1.89123E-04 941 1.27770E-05 942 5.72363E-07
 951 4.85452E-07 925 6.33711E-06 926 4.10550E-07 928 2.58219E-03
 8 8.90358E-03 11 1.03968E-02 26 1.17896E-02 24 3.28650E-03
 28 2.31220E-03 42 2.61937E-04 25 2.74457E-04 884 0.00000E+00
 894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28188E-03
 937 4.55646E-07 939 4.87843E-07 950 1.08070E-08 953 2.69633E-08

962 2.55573E-08 963 8.26626E-10 964 2.49699E-09 965 9.16727E-11
 948 9.442290E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
 967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
 580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
 603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
 608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
 628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
 633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
 648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
 39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
 529 0.00000E+00 555 0.00000E+00 105 2.63858E-04 115 7.74028E-04
 147 0.00000E+00 1 2.17919E-03 402 0.00000E+00 4 0.00000E+00
 13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
 20 0.00000E+00 887 1.50137E-05 897 2.77085E-04 857 1.40819E-06
 817 1.02094E-05 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
 819 0.00000E+00

INCORE
 PREP
 OUTER CORE, GDI1 (OPTIMIZED) EXPANSION COEFFICIENT
 1 1 0 0 0 0 0 3 -20 0 0 0 0 70
 1373.15 0.0 0.0
 85
 1.0
 949 1.65298E-03 940 1.85560E-04 941 8.98239E-06 942 2.81111E-07
 951 3.57843E-07 925 6.30652E-06 926 2.76787E-07 928 2.40805E-03
 8 8.92030E-03 11 1.03968E-02 26 1.17896E-02 24 3.28651E-03
 28 2.31221E-03 42 2.61937E-04 25 2.74458E-04 884 0.00000E+00
 894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28187E-03
 937 3.29710E-07 939 3.14183E-07 950 5.66737E-09 953 9.08185E-09
 962 1.28874E-08 963 2.95563E-10 964 5.87203E-10 965 1.52412E-11
 948 4.78690E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
 967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
 580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
 603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
 608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
 628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
 633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
 648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
 39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
 529 0.00000E+00 555 0.00000E+00 105 2.91455E-04 115 7.74030E-04
 147 0.00000E+00 1 2.17918E-03 402 0.00000E+00 4 0.00000E+00
 13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
 20 0.00000E+00 887 1.04505E-05 897 2.23876E-04 857 9.45563E-07
 817 6.79008E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
 819 0.00000E+00

OTCORE
 PREP
 PLENUM UP&LO, LO Q PU,165%FUEL,
 1 1 0 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0
 6
 1.0
 11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
 42 2.61950E-04 25 2.74470E-04
 PLENUM
 PREP
 PLUGGED PIN, LO Q ,165%FUEL
 1 1 0 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

6
 1.0
 11 1.03970E-02 26 2.96650E-02 24 8.26930E-03 28 5.81800E-03
 42 6.59100E-04 25 6.90610E-04

PLUGGD

PREP

ROD FOLLOWER/GAP,

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

6

1.0
 11 1.99080E-02 26 5.24170E-03 24 1.46120E-03 28 1.02800E-03
 42 1.16460E-04 25 1.22020E-04

FOLLOW

PREP

UPP/RAD'L SH'LD,

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

6

1.0
 11 4.38990E-03 26 4.50900E-02 24 1.25690E-02 28 8.84310E-03
 42 1.00180E-03 25 1.04970E-03

UPPSHD

PREP

LOWER SHIELD,

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

9

1.0
 11 6.36540E-03 26 1.77540E-02 24 4.94900E-03 28 3.48200E-03
 42 3.94460E-04 25 4.13320E-04 6 9.24600E-03 105 7.89610E-03
 115 2.90880E-02

LOWSHD

PREP

ROD ABSORBER HOMOGENEOUS GDI1

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

9

1.0
 11 9.32849E-03 26 1.27380E-02 24 3.55070E-03 28 2.49820E-03
 42 2.83010E-04 25 2.96540E-04 6 8.22779E-03 105 3.04990E-02
 115 2.41210E-03

AHOMOG

PREP

INNER CORE, GDI1 (OPTIMIZED) EXPANSION COEFFICIENT

1 1 0 0 0 0 3 -20 0 0 0 0 70
 1373.15 0.0 0.0

85

1.0
 949 1.40054E-03 940 1.89123E-04 941 1.27770E-05 942 5.72363E-07
 951 4.85452E-07 925 6.33711E-06 926 4.10550E-07 928 2.58219E-03
 8 8.90358E-03 11 1.03968E-02 26 1.17896E-02 24 3.28650E-03
 28 2.31220E-03 42 2.61937E-04 25 2.74457E-04 884 0.00000E+00
 894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28188E-03
 937 4.55646E-07 939 4.87843E-07 950 1.08070E-08 953 2.69633E-08
 962 2.55573E-08 963 8.26626E-10 964 2.49699E-09 965 9.16727E-11
 948 9.44290E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
 967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
 580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00

603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
 608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
 628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
 633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
 648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
 39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
 529 0.00000E+00 555 0.00000E+00 105 2.63858E-04 115 7.74028E-04
 147 0.00000E+00 1 2.17919E-03 402 0.00000E+00 4 0.00000E+00
 13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
 20 0.00000E+00 887 1.50137E-05 897 2.77085E-04 857 1.40819E-06
 817 1.02094E-05 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
 819 0.00000E+00

INCDDD

PREP

OUTER CORE, GDI1 (OPTIMIZED) EXPANSION COEFFICIENT

1 1 0 0 0 0 3 -20 0 0 0 0 70
 1373.15 0.0 0.0

85

1.0

949 1.65298E-03 940 1.85560E-04 941 8.98239E-06 942 2.81111E-07
 951 3.57843E-07 925 6.30652E-06 926 2.76787E-07 928 2.40805E-03
 8 8.92030E-03 11 1.03968E-02 26 1.17896E-02 24 3.28651E-03
 28 2.31221E-03 42 2.61937E-04 25 2.74458E-04 884 0.00000E+00
 894 0.00000E+00 6 2.84691E-04 157 0.00000E+00 40 1.28187E-03
 937 3.29710E-07 939 3.14183E-07 950 5.66737E-09 953 9.08185E-09
 962 1.28874E-08 963 2.95563E-10 964 5.87203E-10 965 1.52412E-11
 948 4.78690E-08 854 0.00000E+00 814 0.00000E+00 966 0.00000E+00
 967 0.00000E+00 968 0.00000E+00 439 0.00000E+00 579 0.00000E+00
 580 0.00000E+00 582 0.00000E+00 584 0.00000E+00 591 0.00000E+00
 603 0.00000E+00 604 0.00000E+00 605 0.00000E+00 606 0.00000E+00
 608 0.00000E+00 600 0.00000E+00 617 0.00000E+00 627 0.00000E+00
 628 0.00000E+00 620 0.00000E+00 622 0.00000E+00 631 0.00000E+00
 633 0.00000E+00 634 0.00000E+00 644 0.00000E+00 646 0.00000E+00
 648 0.00000E+00 82 0.00000E+00 839 0.00000E+00 12 0.00000E+00
 39 0.00000E+00 441 0.00000E+00 442 0.00000E+00 443 0.00000E+00
 529 0.00000E+00 555 0.00000E+00 105 2.91455E-04 115 7.74030E-04
 147 0.00000E+00 1 2.17918E-03 402 0.00000E+00 4 0.00000E+00
 13 0.00000E+00 63 0.00000E+00 73 0.00000E+00 74 0.00000E+00
 20 0.00000E+00 887 1.04505E-05 897 2.23876E-04 857 9.45563E-07
 817 6.79008E-06 889 0.00000E+00 899 0.00000E+00 859 0.00000E+00
 819 0.00000E+00

OTCDDD

PREP

PLENUM UP&LO, LO Q PU, 165%FUEL,

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 1.17900E-02 24 3.28650E-03 28 2.31220E-03
 42 2.61950E-04 25 2.74470E-04

PLEDDD

PREP

PLUGGED PIN, LO Q, 165%FUEL

1 1 0 0 0 0 3 -20 0 0 0 0 70
 703.15 0.0 0.0

6

1.0

11 1.03970E-02 26 2.96650E-02 24 8.26930E-03 28 5.81800E-03
 42 6.59100E-04 25 6.90610E-04

PLUGG 00000120
 FOLLO 00000120
 UPPSH 00000120
 LOWSH 00000120
 AHOMO 00000120

Data file [28] .OPEDID.DATA(PFC3C18X)

condensation JOINT input, 18 groups, normal,
 for 'shutdown margin'

CITATIONPOST	00000010
1	00000020
18	00000030
2 4 6 8 10 13 16 19 22 25 28 31 34 37 40 43 46 70	00000040
PDSIN INCOR SLAR 0 1	00000050
PDSOUT INCOR SLAR	00000060
USERPDS INCOR CITA	00000070
INNER-CORE OPTIMIZED CORE	00000080
PDSIN OTCOR SLAR 0 1	00000090
PDSOUT OTCOR SLAR	00000100
USERPDS OTCOR CITA	00000110
OUTER-CORE OPTIMIZED CORE	00000120
PDSIN PLENU SLAR 0 1	00000130
PDSOUT PLENU SLAR	00000140
USERPDS PLENU CITA	00000150
PLENUM OPTIMIZED CORE	00000160
PDSIN PLUGG SLAR 0 1	00000170
PDSOUT PLUGG SLAR	00000180
USERPDS PLUGG CITA	00000190
STEEL PLUG OPTIMIZED CORE	00000200
PDSIN FOLLO SLAR 0 1	00000210
PDSOUT FOLLO SLAR	00000220
USERPDS FOLLO CITA	00000230
ROD FOLLOWER OPTIMIZED CORE	00000240
PDSIN UPPSH SLAR 0 1	00000250
PDSOUT UPPSH SLAR	00000260
USERPDS UPPSH CITA	00000270
UPPER-SH'D OPTIMIZED CORE	00000280
PDSIN LOWSH SLAR 0 1	00000290
PDSOUT LOWSH SLAR	00000300
USERPDS LOWSH CITA	00000310
RAD'L-SH'D OPTIMIZED CORE	00000320
PDSIN AHOMO SLAR 0 1	00000330
PDSOUT AHOMO SLAR	00000340
USERPDS AHOMO CITA	00000350
ROD ABSORBER OPTIMIZED CORE	00000360

Data file [29] .OPEDID.DATA(PFC3MX)

x-sectio re-formatting JOINT input, normal, for
 'shutdown margin'

CITATIONMICRO	00000100
INCOR SLAROM	00000200
85	00000300
1 949 FU239 239.0530 0.334000E-10	00000400
2 940 FU240 240.0540 0.336 E-10	00000410
3 941 FU241 241.0570 0.337 E-10	00000420
4 942 FU242 242.0580 0.338 E-10	00000420
5 951 AM241 241.0750 0.336 E-10	00000420
6 925 U235 235.0440 0.323 E-10	00000420
7 926 U236 236.0460 0.324 E-10	00000420
8 928 U238 238.0510 0.331 E-10	00000420
9 8 O 15.9994 0.0	00000420
10 11 NA 22.9898 0.0	00000420
11 26 FE 55.8470 0.0	00000420
12 24 CR 51.9961 0.0	00000400
13 28 NI 58.6900 0.0	00000410
14 42 MO 95.9400 0.0	00000420
15 25 MN 54.9380 0.0	00000420
16 884 U-238FP 238.0 0.0	00000420
17 894 FU-239FP 239.0 0.0	00000420
18 6 C 12.011 0.0	00000420
19 157 N15 15.000 0.0	00000420
20 40 ZR 91.224 0.0	
21 937 NF237 237.048 0.330000E-10	
22 939 NF239 239.053 0.331000E-10	
23 950 AM242M 242.060 0.345000E-10	
24 953 AM243 243.061 0.346000E-10	
25 962 CM242 242.059 0.352000E-10	
26 963 CM243 243.060 0.352000E-10	
27 964 CM244 244.063 0.353000E-10	
28 965 CM245 245.066 0.353000E-10	
29 948 FU238 238.050 0.337000E-10	
30 854 U235-FP 235.0 0.0	
31 814 FU241-FP 241.0 0.0	
32 966 CM246 246.0 0.354000E-10	
33 967 CM247 247.0 0.354000E-10	
34 968 CM248 248.0 0.354000E-10	
35 439 TC99 99.0 0.0	
36 579 LA139 139.0 0.0	
37 580 CE140 140.0 0.0	
38 582 CE142 142.0 0.0	
39 584 CE144 144.0 0.0	
40 591 PR141 141.0 0.0	
41 603 ND143 143.0 0.0	
42 604 ND144 144.0 0.0	
43 605 ND145 145.0 0.0	
44 606 ND146 146.0 0.0	
45 608 ND148 148.0 0.0	
46 600 ND150 150.0 0.0	
47 617 PM147 146.915 0.0	
48 627 SM147 147.0 0.0	

00000900

49	628	SM148	148.0	0.0																					
50	620	SM150	150.0	0.0																					
51	622	SM152	152.0	0.0																					
52	631	EU151	150.919	0.0																					
53	633	EU153	152.921	0.0																					
54	634	EU154	154.0	0.0																					
55	644	GD154	154.0	0.0																					
56	646	GD156	155.922	0.0																					
57	648	GD158	157.924	0.0																					
58	82	PB56	207.2	0.0																					
59	839	BI209	209.0	0.0																					
60	12	MG	24.305	0.0																					
61	39	Y	88.9069	0.0																					
62	441	RU101	100.906	0.0																					
63	442	RU102	101.904	0.0																					
64	443	RU103	102.906	0.0																					
65	529	I129	128.905	0.0																					
66	555	CS135	134.906	0.0																					
67	105	B10	10.0129	0.0	00000420																				
68	115	B11	11.0093	0.0	00000420																				
69	147	N14	14.0031	0.0																					
70	1	H	1.0078	0.0																					
71	402	HE	4.0026	0.0																					
72	4	BE	9.0122	0.0																					
73	13	AL27	26.9818	0.0																					
74	63	LI6	6.0151	0.0																					
75	73	LI7	7.0160	0.0																					
76	74	W-NAT	183.8500	0.0																					
77	20	CA	40.0800	0.0																					
78	887	U238-FP	238.0	0.0																					
79	897	PU239-FP	239.0	0.0																					
80	857	U235-FP	235.0	0.0																					
81	817	PU241-FP	241.0	0.0																					
82	889	U238-FP	238.0	0.0																					
83	899	PU239-FP	239.0	0.0																					
84	859	U235-FP	235.0	0.0																					
85	819	PU241-FP	241.0	0.0																					
INCR					85 INNE CORE REGION																				
1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	000000280	
25	26	27	28	29	30	31	32	33	34	35	36	37	38	39	40	41	42	43	44	45	46	47	48		
49	50	51	52	53	54	55	56	57	58	59	60	61	62	63	64	65	66	67	68	69	70	71	72		
73	74	75	76	77	78	79	80	81	82	83	84	85													
OTCOR					85 OUT CORE REGION																				
1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	000000280	
25	26	27	28	29	30	31	32	33	34	35	36	37	38	39	40	41	42	43	44	45	46	47	48		
49	50	51	52	53	54	55	56	57	58	59	60	61	62	63	64	65	66	67	68	69	70	71	72		
73	74	75	76	77	78	79	80	81	82	83	84	85													
PLENU					6 PLENUM																				00000310
10	11	12	13	14	15																			00000280	
PLUGG					6 STEEL PLUG																				00000310
10	11	12	13	14	15																			00000280	
FOLLO					6 ROD FOLLOWER																				00000320
10	11	12	13	14	15																			00000280	
UPPSH					6 UPPER SHIELD																				00000310
10	11	12	13	14	15																			00000280	
LOWSH					9 LOWER SHIELD																				00000310
10	11	12	13	14	15	18	67	68																00000280	
AHOMO					9 ROD ABSORBER																				00000320
10	11	12	13	14	15	18	67	68																00000280	

Data file [30] .OPEDID.DATA(NAMEF#XX)

PERKY package input, material zones, for 'shutdown margin'

- INCR
- OTCOR
- PLENU
- PLUGG
- FOLLO
- UPPSH
- LOWSH
- AHOMO
- INCRD
- OTCDD
- PLEDD
- PLUDD
- FOLDD
- UPPDD
- LOWDD
- AHODD

Data file [31] .OPEDID.DATA(IDNUMXX)

PERKY package input, nuclide IDs, for 'shutdown margin'

- 1 949
- 2 940
- 3 941
- 4 942
- 5 951
- 6 925
- 7 926
- 8 928
- 9 8
- 10 11
- 11 26
- 12 24
- 13 28
- 14 42
- 15 25
- 16 884
- 17 894
- 18 6
- 19 157

20 40
21 937
22 939
23 950
24 953
25 962
26 963
27 964
28 965
29 948
30 854
31 814
32 966
33 967
34 968
35 439
36 579
37 580
38 582
39 584
40 591
41 603
42 604
43 605
44 606
45 608
46 600
47 617
48 627
49 628
50 620
51 622
52 631
53 633
54 634
55 644
56 646
57 548
58 82
59 83
60 12
61 39
62 441
63 442
64 443
65 539
66 555
67 105
68 115
69 147
70 1
71 2
72 4
73 13
74 36
75 37
76 74
77 20
78 887

79 897
80 857
81 817
82 889
83 899
84 859
85 819

Data file [32] .OPEDID.DATA(CHIMG)

PERKY package input, fission spectrum, for 'shutdown margin'

949
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
940
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
941
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
942
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
951
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
925
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
926
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
928
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
937
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
939
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
950

```

2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
953
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
962
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
963
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
964
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
965
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
948
2.75164E-02 1.20304E-01 2.22830E-01 2.34682E-01 1.73023E-01 1.40659E-01
5.36056E-02 1.83258E-02 6.10716E-03 1.99826E-03 6.48725E-04 2.11947E-04
6.87782E-05 2.22527E-05 0.00000E+00 0.00000E+00 0.00000E+00 0.00000E+00
0

```

Data file [33] .OPEDID.DATA(ZONE#XX)

PERKY package input, material zones, for 'shutdown margin'

```

0 /* OUTPUT OPTION */
1 8 INNER CORE
9 32 OUTER CORE
33 41 SHIELD 1
42 45 FOLLOWER
46 48 SHIELD 2
49 52 ABSORBER

```

Data file [34] .OPMOSD.DATA(GDI1MBRN)

MOSES input, burn-up, for '3D burn-up'

GDI1S - LQ PU40/45%,165%F,4*9.75M.,21P ZRH/8P B4C - 18 GROUPS OPTIMIZED
3D BURNUP - SOC 60/50/50/50% EOC 0% STEP 1 CLEAN CORE SOC&B-U
&& MODIFIED - 65/50/50/50 & 55/50/50/50 FOR REFUELLING STEPS 1 & 3 SOC

```

&& CORE VARIANT 7 IMPROVED AXIAL REFLECTORS HETEROGENEOUS RODS
&& -----BORON IFIS=-1----- 1 MESH PER S/A ----- NON-ZERORODTIME
&& SIMPLIFIED CHAIN, NO MA CHAIN, B BURN (WITH F.P. CAPTURE CORRECT)
CONTROL
3*1 && DIFFUSION CAL. ON
3*1 && BURNUP CAL. ON
EDIT
0 4*1 2 19*0 1 4*0 && EDIT PROGRAM SETUP
0 1 3*0 1 3*0 1 20*0 && EDIT FUEL MANAGEMENT DATA
0 0 2 2 1 0 1 0 1 1 0 1 0 1 3 0 0 1 1 0 0 1 3*0 1 0 1 0 0 && DIF CAL
3*0 1 4*0 1 0 2 5*0 9*0 1 4*0 && EDIT BURNUP CAL. DATA

EDITD
&&
&& DETAILED EDIT FOR DIFF. CAL.
50*0.0
&& DETAILED EDIT FOR BURNUP CAL.
50*0.0
&& DETAILED EDIT FOR N.D. CAL.
50*0.0

EOB
&&
&&
&&
LOADING 1
&&
GEOM
&& NSYM NLAY NPLN NAZN NGRP PITCH
2 18 56 15 7 15.81
&& AXIAL MESH BY REGION
10 2 1 2 1 2 3 4 4 3 2 1 1 14 6
&& REGION WIDTH (CM)
50.0 7.0 3.0 5.0 3.0 6.0 9.0 2*12.0 9.0 6.0 3.0 2.0 70.0 33.0
&& === NMES ( NMES=1:CM, NMES=2:MOD.CM)
&& NMES NCAL NITR NIIT NITM NBATCH OUTCR RLOSS DELAY NSMALL
CPARA1 2 0 150 8 1600 0 0 0 0 0 3*0 148.28 1
CPARA2 5.E-5 5.E-4 0.5 1.5 1 0.0 4*0 0.0 0 0.0
EOB
&&
&&
&&
ASSEMBLY AXIAL INFORMATIONS
AXIAL
&& NREGZ
15 6 5 4 3 8*1 4 3 7 && INNER FUEL
50.0 7.0 3.0 5.0 3.0 6.0 9.0 2*12.0 9.0 6.0 3.0 2.0 70.0 33.0
15 6 5 4 3 8*2 4 3 7 && OUTER FUEL
50.0 7.0 3.0 5.0 3.0 6.0 9.0 2*12.0 9.0 6.0 3.0 2.0 70.0 33.0
1 6 && RADIAL SHIELD
230.0
3 8 5 7 && ROD 0%
65.0 132.0 33.0
4 5 8 5 7 && ROD 60%
40.0 61.0 96.0 33.0
4 5 8 5 7 && ROD 50%
35.0 60.0 102.0 33.0
15 6 5 4 3 8*1 4 3 7 && INNER FUELNEW
50.0 7.0 3.0 5.0 3.0 6.0 9.0 2*12.0 9.0 6.0 3.0 2.0 70.0 33.0
15 6 5 4 3 8*2 4 3 7 && OUTER FUELNEW

```

50.0 7.0 3.0 5.0 3.0 6.0 9.0 2*12.0 9.0 6.0 3.0 2.0 70.0 33.0
 4 5 8 5 7 && ROD 65%
 45.0 59.0 93.0 33.0
 4 5 8 5 7 && ROD 55%
 40.0 58.0 99.0 33.0

EOI
 && ASSEMBLY GROUP ASSIGNMENT

ASYGROUP
 1 1 1 1 1 IC1 INNER CORE
 1 2 1 2 2 OC1 OUTER CORE
 1 3 1 4 3 SHD RADIAL SHIELD
 2 4 1 4 -4 R0% ROD 0%
 2 5 1 4 -5 R60 ROD 60%
 2 6 1 4 -6 R50 ROD 50%
 1 7 1 1 1 IC2 INNER CORE2
 1 8 1 2 2 OC2 OUTER CORE2
 2 9 1 4 -9 R65 ROD 65%
 2 10 1 4 -10 R55 ROD 55%

EOI
 && MATERIAL ASSIGNMENT

REGION
 1 1 1 1 1 \$IC1 INNER CORE
 2 2 1 1 1 \$OC1 OUTER CORE
 3 3 1 3 1 \$PLN PLENUM AXIAL
 4 4 1 3 1 \$PLG PLUG AXIAL
 5 5 1 3 1 \$FOL GAP FOLLOWER
 6 6 1 5 1 \$URS UP/RAD SHLD
 7 7 1 3 1 \$LSH LOWER SHIELD
 8 8 1 6 1 \$ABS ROD ABSORBER

EOI
 && CROSS SECTION AND FISSION SPECTRUM

MICROXS
 1 9 33 2 1 2
 1.499E-1 4.330E-1 4.058E-1 1.086E-2 3.839E-4 0.0 0.0

&& NUCLIDE INFORMATION

&& XSNUC

CODE	NAME	IFIS	KIND	BCAL	KLIB	MTW	EFIS	ECAP	DECAY
1	PU239	2	1	0	1	239.05	3.34E-11	0	0
2	PU240	2	2	0	1	240.05	3.36E-11	0	0
3	PU241	2	1	0	1	241.06	3.37E-11	0	1.526E-9
4	PU242	2	3	0	1	242.06	3.38E-11	0	0
5	AM241	2	3	0	1	241.06	3.36E-11	0	0
6	U235	1	3	0	1	235.044	3.23E-11	0	0
7	U236	1	3	0	1	236.05	3.24E-11	0	0
8	U238	1	2	0	1	238.051	3.31E-11	0	0
9	O	0	5	0	1	15.999	0	0	0
10	NA	0	6	0	1	22.990	0	0	0
11	FE	0	6	0	1	55.847	0	0	0
12	CR	0	6	0	1	51.996	0	0	0
13	NI	0	6	0	1	58.70	0	0	0
14	MO	0	6	0	1	95.94	0	0	0
15	MN	0	6	0	1	54.938	0	0	0
18	C	0	7	0	1	12.011	0	0	0
20	ZR	0	6	0	1	91.224	0	0	0
21	NP237	0	3	0	1	237.048	3.30E-11	0	0

22	NP239	0	3	0	1	239.053	3.31E-11	0	3.406E-6
23	AM242M	0	3	0	1	242.060	3.45E-11	0	0
24	AM243	0	3	0	1	243.061	3.46E-11	0	0
25	CM242	0	3	0	1	242.059	3.52E-11	0	4.922E-8
26	CM243	0	3	0	1	243.060	3.52E-11	0	0
27	CM244	0	3	0	1	244.063	3.53E-11	0	0
28	CM245	0	3	0	1	245.066	3.53E-11	0	0
29	PU238	2	2	0	1	238.05	3.33E-11	0	0
67	B-10	-1	7	0	1	10.0123	0	0	0
68	B-11	-1	7	0	1	11.0093	0	0	0
70	H	0	6	0	1	1.0078	0	0	0
78	U238-FP	-1	4	0	1	238.051	0	0	0
79	PU239-FP	-1	4	0	1	239.055	0	0	0
80	U235-FP	-1	4	0	1	235.044	0	0	0
81	PU241-FP	-1	4	0	1	241.057	0	0	0

REACT1

6	7	2	1.0
8	1	2	1.0
1	2	2	1.0
2	3	2	1.0
3	4	2	1.0
29	1	2	1.0
3	5	1	1.0
67	68	2	1.0
78	78	1	1.0
79	79	1	1.0
80	80	1	1.0
81	81	1	1.0
0			

YIELD

78	8	1.0
79	1	1.0
79	2	1.0
79	29	1.0
80	6	1.0
80	7	1.0
81	3	1.0
81	4	1.0
81	5	1.0

&& ATOM NUMBER DENSITY

&& ATDEN

&& FROM GD11 PENCIL OUTPUT/CITDENS INPUT

&& ZONE 1

1.69700E-03	1.07870E-04	0.0	0.0	8.17870E-06
0.0	2.68370E-03	8.90360E-03	1.03968E-02	1.17896E-02
2.31220E-03	2.61937E-04	2.74457E-04	2.8469E-04	1.28190E-03
0.0	0.0	0.0	0.0	0.0
0.0	0.0	3.6474E-04	7.7403E-04	2.17920E-03
0.0	0.0	0.0	0.0	0.0

&& ZONE 2

1.90280E-03	1.20950E-04	0.0	0.0	7.53940E-06
0.0	2.47390E-03	8.92030E-03	1.03968E-02	1.17896E-02
2.31220E-03	2.61937E-04	2.74457E-04	2.8469E-04	1.28190E-03
0.0	0.0	0.0	0.0	0.0
0.0	0.0	3.6474E-04	7.7403E-04	2.17920E-03
0.0	0.0	0.0	0.0	0.0

&& ZONE 3

96	96	8
98	98	8
104	104	8
106	106	8
108	108	8
113	113	8
122	122	8
124	124	8
133	133	8
138	138	8
142	142	8
150	150	8
154	154	8
159	159	8
161	161	8
170	170	8
172	172	8
174	174	8
183	183	8
190	190	8
192	192	8
204	204	8
206	206	8
8	8	9
11	11	9
35	35	6
41	41	6
74	74	6
77	77	6
80	80	6
83	83	6
86	86	6
89	89	6
137	137	6
143	143	6
149	149	6
155	155	6
0	0	0
4THCORE		
3	3	7
10	10	7
13	13	7
23	23	7
25	25	7
27	27	7
29	29	7
31	31	7
45	45	7
47	47	7
49	49	7
51	51	7
53	53	7
55	55	7
57	57	7
78	78	7
81	81	7
87	87	7
90	90	7
93	93	8

101	101	8
103	103	8
111	111	8
116	116	8
118	118	8
120	120	8
127	127	8
129	129	8
131	131	8
135	135	8
145	145	8
147	147	8
157	157	8
162	162	8
164	164	8
166	166	8
168	168	8
175	175	8
177	177	8
179	179	8
181	181	8
189	189	8
203	203	8
8	8	5
11	11	5
35	35	6
41	41	6
74	74	6
77	77	6
80	80	6
83	83	6
86	86	6
89	89	6
137	137	6
143	143	6
149	149	6
155	155	6
0	0	0
5THCORE		
5	5	7
7	7	7
15	15	7
17	17	7
19	19	7
21	21	7
33	33	7
37	37	7
39	39	7
43	43	7
59	59	7
61	61	7
63	63	7
65	65	7
67	67	7
69	69	7
71	71	7
73	73	7
75	75	7
84	84	7

95	95	8
97	97	8
99	99	8
105	105	8
107	107	8
109	109	8
112	112	8
114	114	8
121	121	8
123	123	8
125	125	8
132	132	8
139	139	8
141	141	8
151	151	8
153	153	8
158	158	8
160	160	8
169	169	8
171	171	8
173	173	8
182	182	8
191	191	8
205	205	8
8	8	10
11	11	10
35	35	6
41	41	6
74	74	6
77	77	6
80	80	6
83	83	6
86	86	6
89	89	6
137	137	6
143	143	6
149	149	6
155	155	6
0	0	0
6THCORE		
2	2	7
9	9	7
12	12	7
22	22	7
24	24	7
26	26	7
28	28	7
30	30	7
44	44	7
46	46	7
48	48	7
50	50	7
52	52	7
54	54	7
56	56	7
76	76	7
79	79	7
85	85	7
88	88	7

92	92	8					
100	100	8					
102	102	8					
110	110	8					
115	115	8					
117	117	8					
119	119	8					
126	126	8					
128	128	8					
130	130	8					
134	134	8					
136	136	8					
144	144	8					
146	146	8					
148	148	8					
156	156	8					
163	163	8					
165	165	8					
167	167	8					
176	176	8					
178	178	8					
180	180	8					
193	193	8					
207	207	8					
8	8	5					
11	11	5					
35	35	6					
41	41	6					
74	74	6					
77	77	6					
80	80	6					
83	83	6					
86	86	6					
89	89	6					
137	137	6					
143	143	6					
149	149	6					
155	155	6					
0	0	0					
EOI							
PLANT							
520.0 380							
CRPOS							
&&	IC1	OC1	RS1	BCR	30%	PCR	
	0.0	0.0	0.0	0.0	0.0	0.0	34*0.0
EOB							
STEP2 - CYCLE 1, MOC & B-U TO EOC							
RODS 0%							
CONTROL							
3*1							
3*1							
EOB							
LOADING 2							
EOB							
CRPOS							
&&	IC1	OC1	RS1	BCR	30%	PCR	
	0.0	0.0	0.0	0.0	0.0	0.0	34*0.0
EOB							
STEP 3 - CYCLE 1, EOC (NO BURN)							

```

RODS 0%
CONTROL
  3*1
  3*0
EOB
LOADING 0
CPARA1 2 0 150 3 1600 0 0 0 0 0 3*0 0.0 0
EOB
PLANT
520.0 380
EOB
STEP 4 - CYCLE 2, REFUEL STEP 1 (3RDCORE)
RODS REVERT TO 65/50/50/50% SOC & B-U TO MOC
CONTROL
  3*1
  3*1
EOB
LOADING 3
CPARA1 2 0 150 3 1600 0 0 0 0 0 3*0 148.28 1
EOB
PLANT
520.0 380
CRPOS
  && IC1 OC1 RS1 BCR 30% PCR
  0.0 0.0 0.0 0.0 0.0 0.0 34*0.0
EOB
STEPS - CYCLE 2, MOC & B-U TO EOC
RODS 0%
CONTROL
  3*1
  3*1
EOB
LOADING 2
EOB
CRPOS
  && IC1 OC1 RS1 BCR 30% PCR
  0.0 0.0 0.0 0.0 0.0 0.0 34*0.0
EOB
STEP 6 - CYCLE 2, EOC (NO BURN)
RODS 0%
CONTROL
  3*1
  3*0
EOB
LOADING 0
CPARA1 2 0 150 3 1600 0 0 0 0 0 3*0 0.0 0
EOB
PLANT
520.0 380
EOB
STEP 7 - CYCLE 3, REFUEL STEP 2 (4THCORE)
RODS REVERT TO 60/50/50/50% SOC & B-U TO MOC
CONTROL
  3*1
  3*1
EOB
LOADING 4
CPARA1 2 0 150 3 1600 0 0 0 0 0 3*0 148.28 1
EOB

```

```

PLANT
520.0 380
CRPOS
  && IC1 OC1 RS1 BCR 30% PCR
  0.0 0.0 0.0 0.0 0.0 0.0 34*0.0
EOB
.....
REPETITIVE DATA REMOVED
.....
STEP 27 - CYCLE 9, EOC (NO BURN)
RODS 0%
CONTROL
  3*1
  3*0
EOB
LOADING 0
CPARA1 2 0 150 3 1600 0 0 0 0 0 3*0 0.0 0
EOB
PLANT
520.0 380
EOB

```

.ARCHTAPE.CNTL(REST30)

ARCHIVE restoration of JCL & data files,
from cartridge SNHBU3

```

//POCOB17X JOB ( ),CLASS=A,MSGCLASS=X,MSGLEVEL=(1,1),
// NOTIFY=POCOB17
//*
//* MT DASD s hi
//*
//*ROUTE PRINT X
//*
//STEP0056 EXEC PGM=JSDGENER
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=DELAY.DATA,DISP=(OLD,PASS),LABEL=(56,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DCB=4
//SYSUT2 DD DSN=POCOB17.DELAY.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(1,9),RLSE)
/*
//STEP0057 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPCITD.DATA,DISP=(OLD,PASS),LABEL=(57,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DCB=4
//SYSUT2 DD DSN=POCOB17.OPCITD.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(4,2,50),RLSE)
/*
//STEP0058 EXEC PGM=JSECOPY

```

```

//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPDEND.DATA,DISP=(OLD,PASS),LABEL=(58,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.OPDEND.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(20,5,50),RLSE)
/*
//STEP0059 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPEDID.DATA,DISP=(OLD,PASS),LABEL=(59,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.OPEDID.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(20,1,50),RLSE)
/*
//STEP0060 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPFORT.DATA,DISP=(OLD,PASS),LABEL=(60,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.OPFORT.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(22,4,50),RLSE)
/*
//STEP0061 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPJCL.CNTL,DISP=(OLD,PASS),LABEL=(61,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.OPJCL.CNTL,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(36,0,50),RLSE)
/*
//STEP0062 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=OPMOSD.DATA,DISP=(OLD,PASS),LABEL=(62,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.OPMOSD.DATA,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(59,2,50),RLSE)
/*
//STEP0063 EXEC PGM=JSDGENER
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=TSS.CLIST,DISP=(OLD,PASS),LABEL=(63,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.TSS.CLIST,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(1,0),RLSE)
/*
//STEP0064 EXEC PGM=JSECOPY
//SYSPRINT DD SYSOUT=*
//SYSIN DD DUMMY
//SYSUT1 DD DSN=TSSMAC.CLIST,DISP=(OLD,PASS),LABEL=(64,SL),
// VOL=SER=SNHBU3,UNIT=CMT0,DCB=DEN=4
//SYSUT2 DD DSN=POC0B17.TSSMAC.CLIST,
// DISP=(,CATLG),UNIT=SYSDA,SPACE=(TRK,(11,0,50),RLSE)
/*
//

```

REFERENCES

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- 2 T.B.Fowler et al Nuclear reactor core analysis code: CITATION ORNL-TM-2496, Rev 2
- 3 S.N.Hunter EDSLA: an input editing program for SLAROM (to be published)
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- 5 S.Iijima, H.Yoshida, H.Sakuragi Calculation program for fast reactor design, 2 (Multi-dimensional perturbation theory code based on diffusion approximation: PERKY) JAERI-M 6993
- 6 D.Saphier et al Evaluated delayed neutron spectra and their importance in reactor calculations NSE 62 770
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- 10 H.Hayashi et al Progress report of the design study on a large scale reactor - design study on a 600 MWe class plant PNC TN9410 92-137